

10/513699

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NEWS 1 Web Page for STN Seminar Schedule - N. America
NEWS 2 AUG 10 Time limit for inactive STN sessions doubles to 40
minutes
NEWS 3 AUG 18 COMPENDEX indexing changed for the Corporate Source
(CS) field
NEWS 4 AUG 24 ENCOMPLIT/ENCOMPLIT2 reloaded and enhanced
NEWS 5 AUG 24 CA/CAPLUS enhanced with legal status information for
U.S. patents
NEWS 6 SEP 09 50 Millionth Unique Chemical Substance Recorded in
CAS REGISTRY
NEWS 7 SEP 11 WPIDS, WPINDEX, and WPIX now include Japanese FTERM
thesaurus
NEWS 8 OCT 21 Derwent World Patents Index Coverage of Indian and
Taiwanese Content Expanded
NEWS 9 OCT 21 Derwent World Patents Index enhanced with human
translated claims for Chinese Applications and
Utility Models
NEWS 10 OCT 27 Free display of legal status information in CA/CAPLUS,
USPATFULL, and USPAT2 in the month of November.
NEWS 11 NOV 23 Addition of SCAN format to selected STN databases
NEWS 12 NOV 23 Annual Reload of IFI Databases

NEWS EXPRESS MAY 26 09 CURRENT WINDOWS VERSION IS V8.4,
AND CURRENT DISCOVER FILE IS DATED 06 APRIL 2009.

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* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 13:44:46 ON 23 NOV 2009

<12/04/2007>

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=> file reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.22

0.22

FILE 'REGISTRY' ENTERED AT 13:44:55 ON 23 NOV 2009

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STRUCTURE FILE UPDATES: 22 NOV 2009 HIGHEST RN 1193309-59-9

DICTIONARY FILE UPDATES: 22 NOV 2009 HIGHEST RN 1193309-59-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 26, 2009.

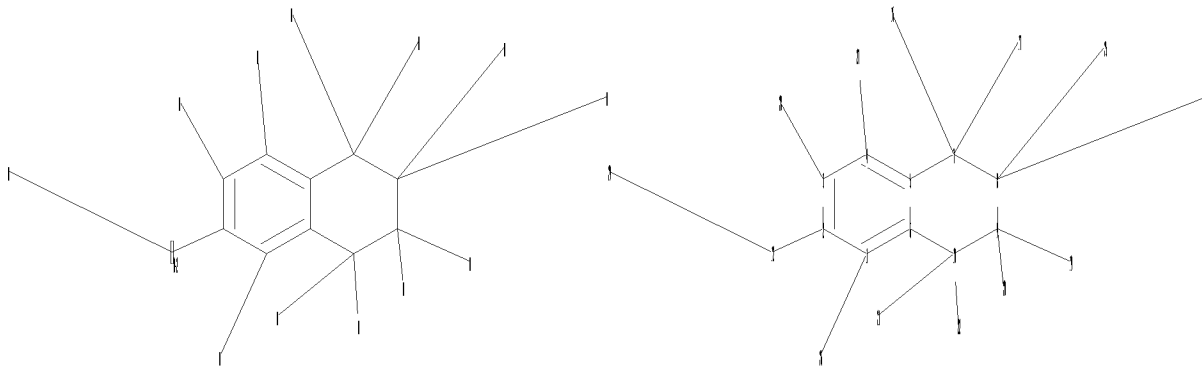
Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10598262hydrogen.str



chain nodes :

12 16 17 19 20 21 22 24 25 26 27 28 29

ring nodes :

1 2 3 4 5 6 7 8 9 10

chain bonds :

1-26 2-12 3-28 4-27 7-16 7-17 8-24 8-25 9-19 9-20 10-21 10-22 12-29

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-10 7-8 8-9 9-10

exact/norm bonds :

<12/04/2007>

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10/513699

5-7 6-10 7-8 8-9 9-10 12-29

exact bonds :

1-26 2-12 3-28 4-27 7-16 7-17 8-24 8-25 9-19 9-20 10-21 10-22

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6

G1:H,N

G2:C,H

G3:C,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom

12:CLASS 16:CLASS 17:CLASS 19:CLASS 20:CLASS 21:CLASS 22:CLASS 24:CLASS

25:CLASS 26:CLASS 27:CLASS 28:CLASS 29:CLASS

L1 STRUCTURE UPLOADED

=> s l1 sss

SAMPLE SEARCH INITIATED 13:45:12 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 188474 TO ITERATE

1.1% PROCESSED 2000 ITERATIONS

3 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
BATCH **INCOMPLETE**

PROJECTED ITERATIONS: 3743870 TO 3795090

PROJECTED ANSWERS: 4646 TO 6662

L2 3 SEA SSS SAM L1

=> s l1 full

THE ESTIMATED SEARCH COST FOR FILE 'REGISTRY' IS 185.40 U.S. DOLLARS

DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N or END:y

FULL SEARCH INITIATED 13:45:19 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 3777563 TO ITERATE

51.7% PROCESSED 1954366 ITERATIONS

4603 ANSWERS

52.9% PROCESSED 2000000 ITERATIONS

4603 ANSWERS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.20

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
BATCH **INCOMPLETE**

PROJECTED ITERATIONS: 3777563 TO 3777563

PROJECTED ANSWERS: 8415 TO 8973

L3 4603 SEA SSS FUL L1

<12/04/2007>

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=> file caplus
COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
185.88	186.10

FULL ESTIMATED COST

FILE 'CAPLUS' ENTERED AT 13:45:43 ON 23 NOV 2009
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FILE COVERS 1907 - 23 Nov 2009 VOL 151 ISS 22
FILE LAST UPDATED: 22 Nov 2009 (20091122/ED)
REVISED CLASS FIELDS (/NCL) LAST RELOADED: Aug 2009
USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Aug 2009

Caplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2009.

CAS Information Use Policies apply and are available at:

<http://www.cas.org/legal/infopolicy.html>

This file contains CAS Registry Numbers for easy and accurate substance identification.

During November, try the new LSUS format of legal status information in the CA/Caplus family databases for free! Complete details on the number of free displays and other databases participating in this offer appear in NEWS 10.

=> s l3 full
L4 116 L3

=> s py<2005
SYSTEM LIMITS EXCEEDED - SEARCH ENDED
The search profile you entered was too complex or gave too many answers. Simplify or subdivide the query and try again. If you have exceeded the answer limit, enter DELETE HISTORY at an arrow prompt (=>) to remove all previous answers sets and begin at L1. Use the SAVE command to store any important profiles or answer sets before using DELETE HISTORY.

=> s l4 py<2005
MISSING OPERATOR L4 PY<2005
The search profile that was entered contains terms or nested terms that are not separated by a logical operator.

<12/04/2007>

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=> s l4 and py<2005

25152347 PY<2005

L5 47 L4 AND PY<2005

=> d ibib abs hitstr tot

THE ESTIMATED COST FOR THIS REQUEST IS 265.08 U.S. DOLLARS

DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N:y

<12/04/2007>

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L5 ANSWER 1 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2005:570893 CAPLUS

DOCUMENT NUMBER: 143:97353

TITLE: Preparation of pyrazolo[3,4-b]pyridines as phosphodiesterase, especially PDE4B, inhibitors for treatment of inflammatory and/or allergic diseases

INVENTOR(S): Allen, David George; Coe, Diane Mary; Cook, Caroline Mary; Dowle, Michael Dennis; Edlin, Christopher David; Hamblin, Julie Nicole; Johnson, Martin Redpath; Jones, Paul Spencer; Lindvall, Mika Kristian; Mitchell, Charlotte Jane; Redgrave, Alison Judith; Robinson, John Edward; Trivedi, Naimisha

PATENT ASSIGNEE(S): Glaxo Group Limited, UK

SOURCE: PCT Int. Appl., 295 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 3

PATENT INFORMATION:

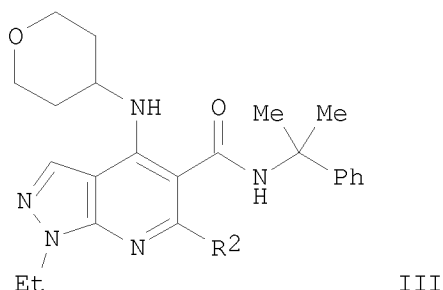
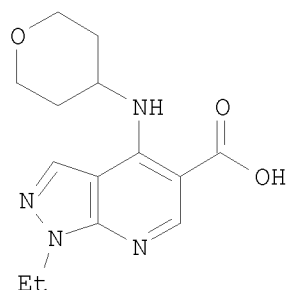
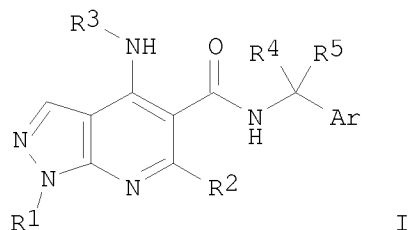
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WO 2004056823	A1	20040708	WO 2003-EP14867	20031219 <--
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AU 2004299277	A1	20050630	AU 2004-299277	20041217
CA 2557004	A1	20050630	CA 2004-2557004	20041217
EP 1737857	A1	20070103	EP 2004-804089	20041217
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CN 1914205	A	20070214	CN 2004-80041657	20041217
JP 2007514704	T	20070607	JP 2006-544380	20041217
US 20070111995	A1	20070517	US 2006-596561	20060616
NO 2006003340	A	20060912	NO 2006-3340	20060718
US 20080132536	A1	20080605	US 2008-22372	20080130
PRIORITY APPLN. INFO.:			WO 2003-EP14867	A 20031219
			GB 2004-5899	A 20040316

<12/04/2007>

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GB 2004-5936	A 20040316
GB 2004-6754	A 20040325
GB 2002-30045	A 20021223
GB 2002-30165	A 20021224
GB 2003-7998	A 20030407
WO 2004-EP14490	W 20041217
US 2006-596561	A1 20060616

OTHER SOURCE(S): CASREACT 143:97353; MARPAT 143:97353
GI



- AB Title compds. I [wherein Ar = (un)substituted Ph, naphthyl, indanyl, pyridinyl, etc. with provisos; R1 = fluoro/alkyl, CH₂CH₂OH; R2 = H, Me, fluoroalkyl; R3 = (un)substituted cycloalkyl, monounsaturated cycloalkenyl, bicyclic, heterocyclic; R4 = H, Me, Et, Pr, i-Pr, fluoroalkyl, cyclopropyl, etc.; R5 = H, fluoro/alkyl, (un)substituted cyclo/alkyl, Ph, etc.; provided that at least one of R4 and R5 is not H; and salts thereof] were prepared as selective phosphodiesterase 4 (PDE4), especially PDE4B, inhibitors. The invention also provides for the use of I for the treatment and/or prophylaxis of an inflammatory and/or allergic disease, such as chronic obstructive pulmonary disease (COPD), asthma, rheumatoid arthritis, allergic rhinitis or atopic dermatitis. Eight biol. methods are given. For example, coupling of acid II with 2-phenyl-2-propanamine gave title compound III. Selected I inhibited PDE4B with IC₅₀ in the range of. I are in particle size-reduced form (DC₅₀ value of about 0.5 to about 10 μ) when used in inhalant compns.
- IT 856559-79-0P, 1-Ethyl-N-[1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]-4-[(tetrahydro-2H-pyran-4-yl)amino]-1H-pyrazolo[3,4-b]pyridine-5-carboxamide 856560-45-7P, 4-(Cyclohexylamino)-1-ethyl-N-[1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]-1H-pyrazolo[3,4-b]pyridine-5-carboxamide 856561-00-7P, 1-Ethyl-4-[(4-oxocyclohexyl)amino]-N-[1-(5,6,7,8-tetrahydro-2-

10/513699

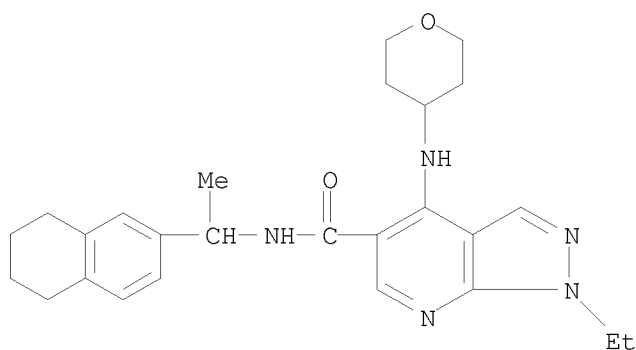
naphthalenyl)ethyl]-1H-pyrazolo[3,4-b]pyridine-5-carboxamide
856561-03-0P, 1-Ethyl-4-[[4-(hydroxyimino)cyclohexyl]amino]-N-[1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]-1H-pyrazolo[3,4-b]pyridine-5-carboxamide

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(PDE4B inhibitor; preparation of pyrazolo[3,4-b]pyridines as PDE4 inhibitors for treatment of inflammatory and/or allergic diseases)

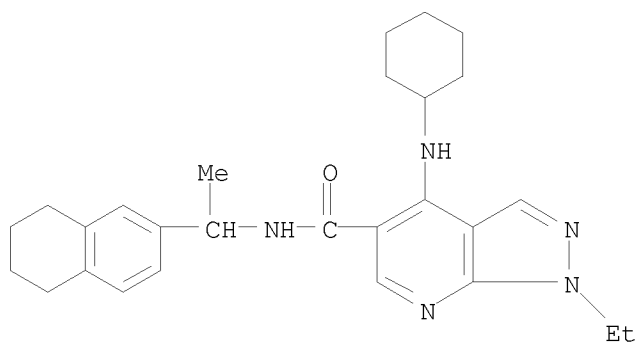
RN 856559-79-0 CAPLUS

CN 1H-Pyrazolo[3,4-b]pyridine-5-carboxamide,
1-ethyl-N-[1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]-4-[(tetrahydro-2H-pyran-4-yl)amino]- (CA INDEX NAME)



RN 856560-45-7 CAPLUS

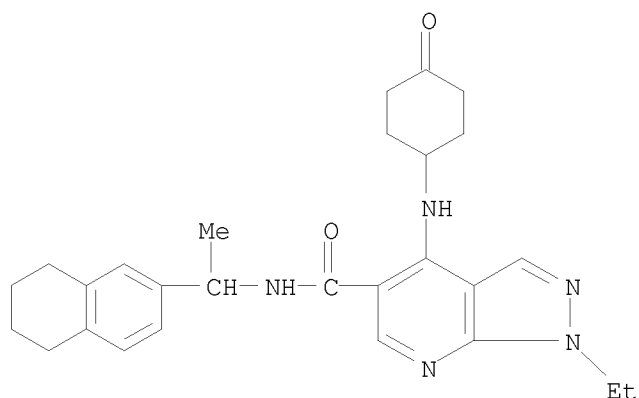
CN 1H-Pyrazolo[3,4-b]pyridine-5-carboxamide,
4-(cyclohexylamino)-1-ethyl-N-[1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]-
(CA INDEX NAME)



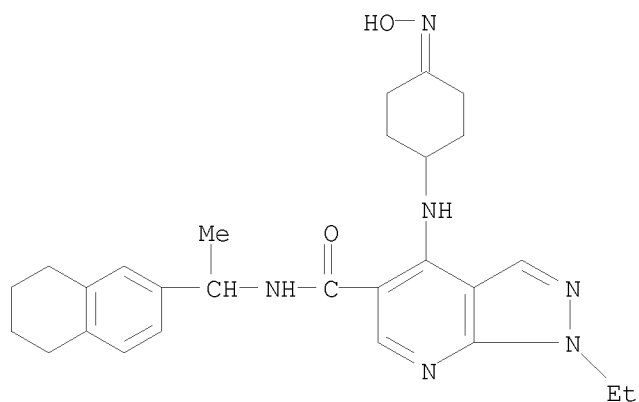
RN 856561-00-7 CAPLUS

CN 1H-Pyrazolo[3,4-b]pyridine-5-carboxamide,
1-ethyl-4-[(4-oxocyclohexyl)amino]-N-[1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]- (CA INDEX NAME)

10/513699



RN 856561-03-0 CAPLUS
CN 1H-Pyrazolo[3,4-b]pyridine-5-carboxamide,
1-ethyl-4-[[4-(hydroxyimino)cyclohexyl]amino]-N-[1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]- (CA INDEX NAME)



OS.CITING REF COUNT: 4 THERE ARE 4 CAPLUS RECORDS THAT CITE THIS RECORD
(4 CITINGS)
REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

10/513699

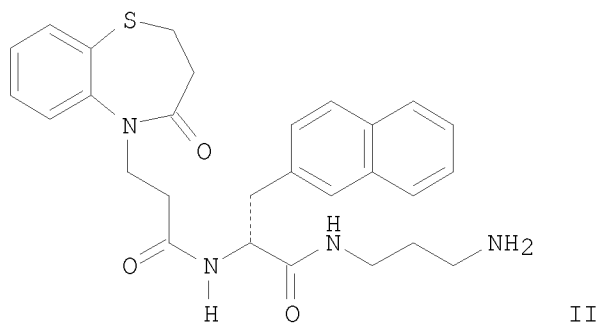
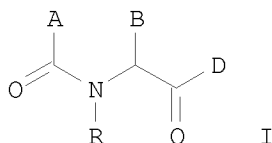
L5 ANSWER 2 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 2005:207843 CAPLUS
DOCUMENT NUMBER: 142:261793
TITLE: Preparation of N-acylated lipophilic amino acid
derivatives having growth hormone releasing activity
INVENTOR(S): Funamizu, Hidenori; Ishiyama, Nobuo; Ikegami, Satoru;
Okuno, Tadashi; Inoguchi, Kiyoshi; Huang, Ping; Loew,
Gilda
PATENT ASSIGNEE(S): Kaken Pharmaceutical Co., Ltd., Japan; Molecular
Research Institute
SOURCE: U.S., 49 pp., Cont.-in-part of U.S. Ser. No. 916,575,
abandoned.
CODEN: USXXAM
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 2
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 6864250	B1	20050308	US 2000-485845	20000426
WO 9909991	A1	19990304	WO 1998-US17232	19980820 <--
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RW:	GH, GM, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG			
US 20050059699	A1	20050317	US 2004-962598	20041013
US 7064121	B2	20060620		
US 20060142264	A1	20060629	US 2006-359616	20060223
US 7279573	B2	20071009		
US 20080027038	A1	20080131	US 2007-892063	20070820
PRIORITY APPLN. INFO.:			US 1997-916575	B2 19970822
			WO 1998-US17232	W 19980820
			US 2000-485845	A3 20000426
			US 2004-962598	A3 20041013
			US 2006-359616	A3 20060223

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): CASREACT 142:261793; MARPAT 142:261793

GI



AB The invention relates to amino acid derivs. I [A is a lipophilic group including an aliphatic bridging group, B is a lipophilic group, D is a group having at least one (un)substituted amino group, R is H, alkyl or cycloalkyl] and their pharmaceutically acceptable salts and individual isomers which have growth hormone releasing activity in humans or animals and are useful, e.g., in treating osteoporosis, bone fractures, wounds or burns. Thus, a 2-step synthesis afforded amide II.HCl, which showed growth hormone releasing activity $< 10^{-8}$ M.

IT 1042394-85-3

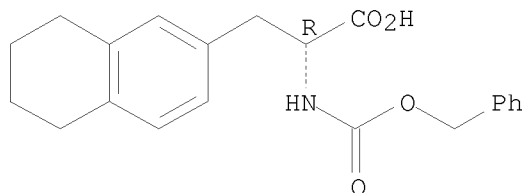
RL: RCT (Reactant); RACT (Reactant or reagent)

(preparation of N-acylated lipophilic amino acid derivs. having growth hormone releasing activity)

RN 1042394-85-3 CAPLUS

CN 2-Naphthalenepropanoic acid, 5,6,7,8-tetrahydro- α -
[[(phenylmethoxy)carbonyl]amino]-, (α R)- (CA INDEX NAME)

Absolute stereochemistry.



OS.CITING REF COUNT: 1 THERE ARE 1 CAPLUS RECORDS THAT CITE THIS RECORD (1 CITINGS)

REFERENCE COUNT: 9 THERE ARE 9 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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L5 ANSWER 3 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2004:333686 CAPLUS

DOCUMENT NUMBER: 140:357056

TITLE: Preparation of novel propargyl ether derivatives for
controlling phytopathogenic microorganisms

INVENTOR(S): Lamberth, Clemens; Zeller, Martin

PATENT ASSIGNEE(S): Syngenta Participations Ag, Switz.

SOURCE: PCT Int. Appl., 57 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

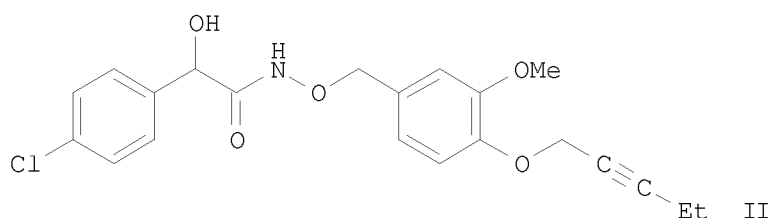
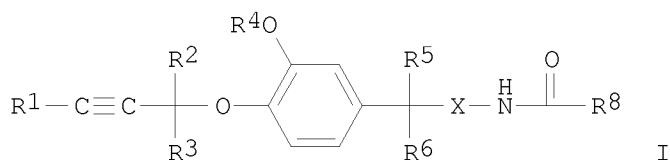
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RW:	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
CA 2498940	A1	20040422	CA 2003-2498940	20031009 <--
AU 2003293609	A1	20040504	AU 2003-293609	20031009 <--
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EP 1549609	B1	20070124		
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US 20060167316	A1	20060727	US 2005-528668	20050322
US 7189873	B2	20070313		
IN 2005CN00561	A	20070622	IN 2005-CN561	20050406
MX 2005003707	A	20050617	MX 2005-3707	20050407
PRIORITY APPLN. INFO.:			GB 2002-23665	A 20021010
			WO 2003-EP11218	W 20031009

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 140:357056

GI



AB The title compds. [I; R1= H, (un)substituted alkyl, cycloalkyl, aryl; R2, R3, R5-R7 = H, alkyl; R4 = (un)substituted alkyl; X = O, NR7; R8 = CR9R10OR11, CR12R13NHSO2R14 (wherein R9 = (un)substituted (hetero)aryl; R10, R11 = H, (un)substituted alkyl, alkenyl, alkynyl; R12 = (un)substituted alkyl, cycloalkyl, aryl, heteroaryl; R13 = H, (un)substituted alkyl, alkenyl or alkynyl; R14 = (un)substituted alkyl, NH2)] which possess plant protecting properties and may advantageously be employed in agricultural practice for controlling or preventing the infestation of plants by phytopathogenic microorganisms, especially fungi, were prepared E.g, a multi-step synthesis of II, starting from 4-hydroxymethyl-2-methoxyphenol and MeCH2C.tplbond.CCH2OH, was given. Representative compds. I showed at least 80% inhibition of fungal infestation in 3 biol. tests.

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	1055764-29-8	1055764-30-1	1055764-31-2
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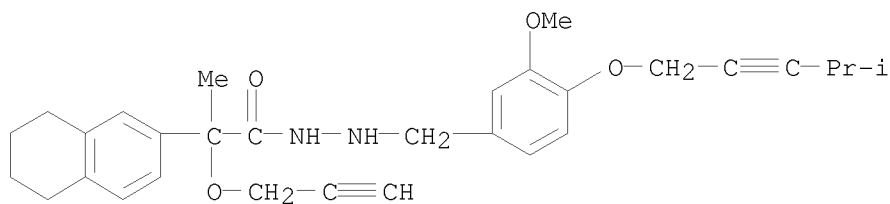
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RL: PRPH (Prophetic)

(Preparation of novel propargyl ether derivatives for controlling
phytopathogenic microorganisms)

RN 1055763-20-6 CAPLUS

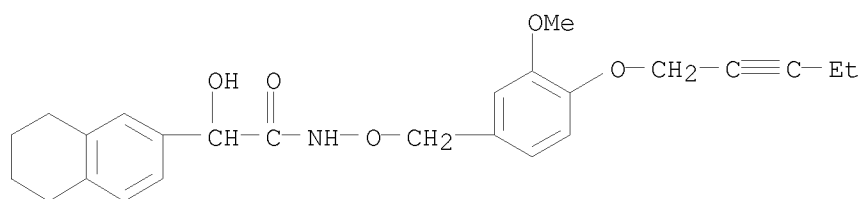
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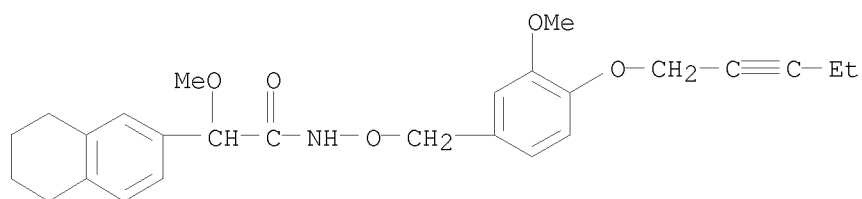
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -hydroxy-N-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methoxy]- (CA INDEX NAME)

10/513699



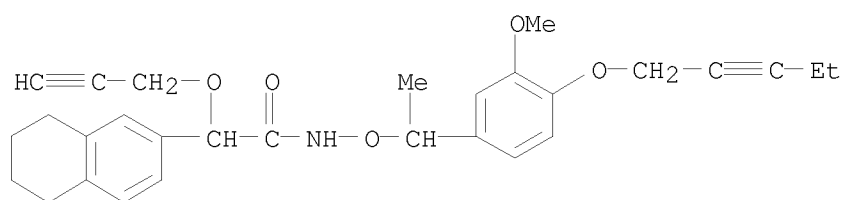
RN 1055764-25-4 CAPLUS

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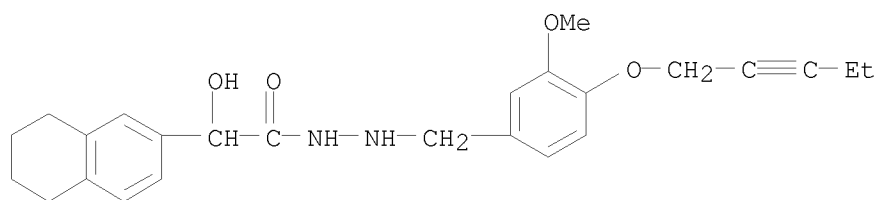
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CN INDEX NAME NOT YET ASSIGNED



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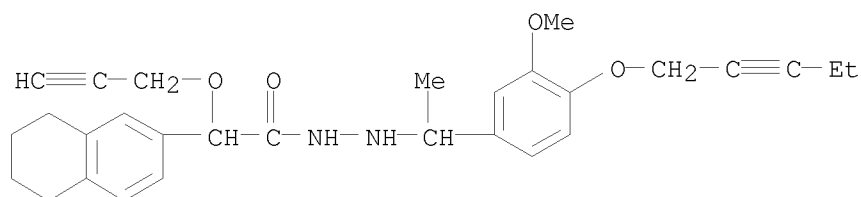
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-, 2-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055764-28-7 CAPLUS

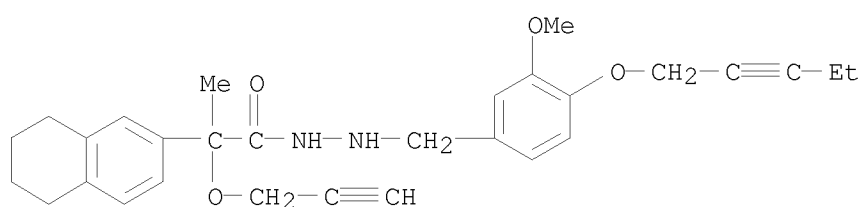
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[1-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]hydrazide (CA INDEX NAME)

10/513699



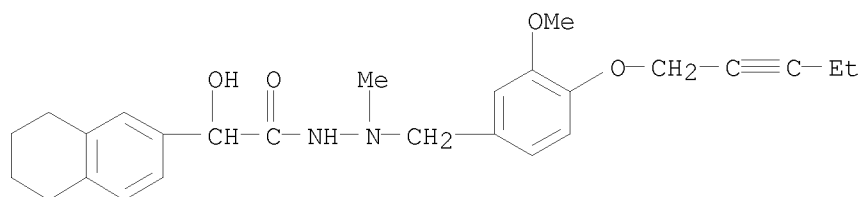
RN 1055764-29-8 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methyl]hydrazide (CA INDEX NAME)



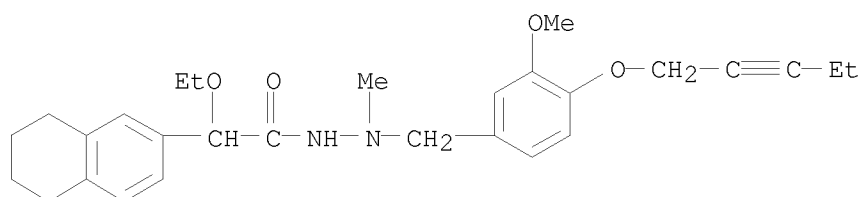
RN 1055764-30-1 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-, 2-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



RN 1055764-31-2 CAPLUS

CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-, 2-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)

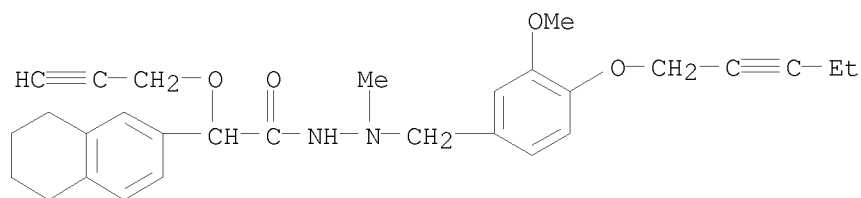


RN 1055764-32-3 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methyl]hydrazide (CA INDEX NAME)

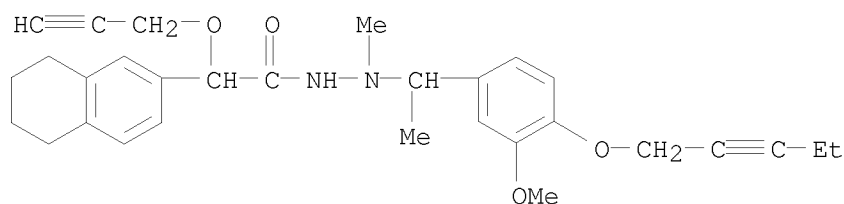
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2-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methyl]-2-methylhydrazide (CA
INDEX NAME)



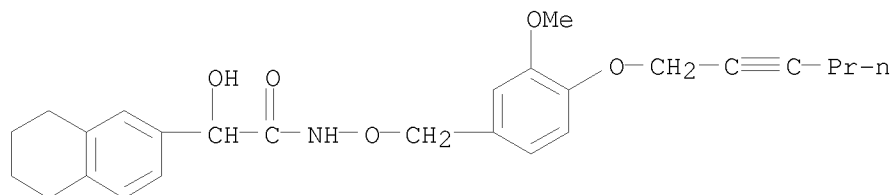
RN 1055764-33-4 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[1-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]-2-methylhydrazide (CA
INDEX NAME)



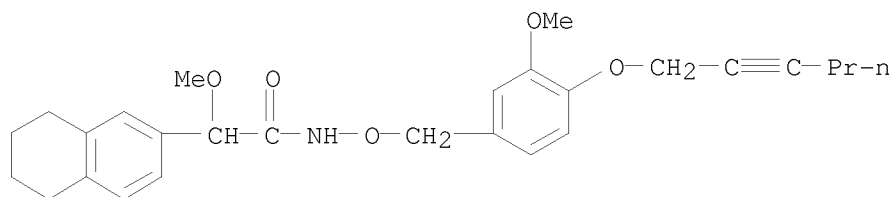
RN 1055765-35-9 CAPLUS

CN 2-Naphthaleneacetamide, N-[[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]methoxy]-
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RN 1055765-37-1 CAPLUS

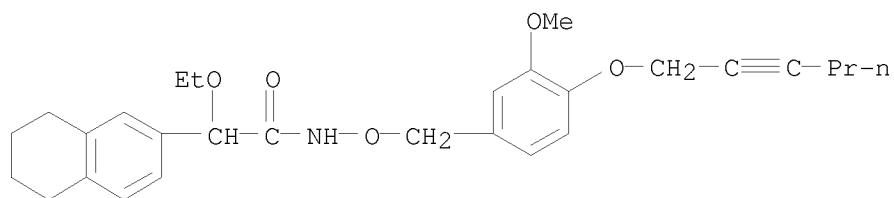
CN 2-Naphthaleneacetamide, N-[[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]methoxy]-
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RN 1055765-38-2 CAPLUS

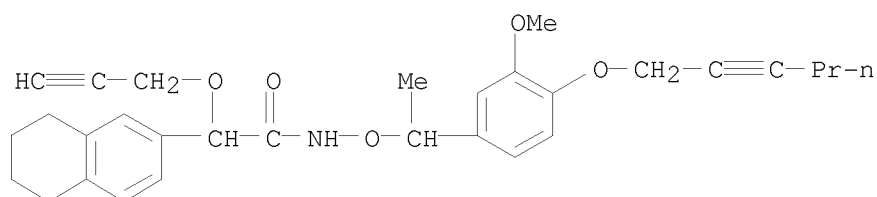
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CN 2-Naphthaleneacetamide, α -ethoxy-N-[[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro- (CA INDEX NAME)



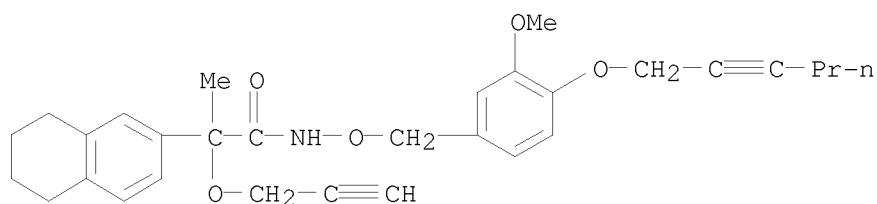
RN 1055765-39-3 CAPLUS

CN 2-Naphthaleneacetamide, N-[1-[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]ethoxy]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



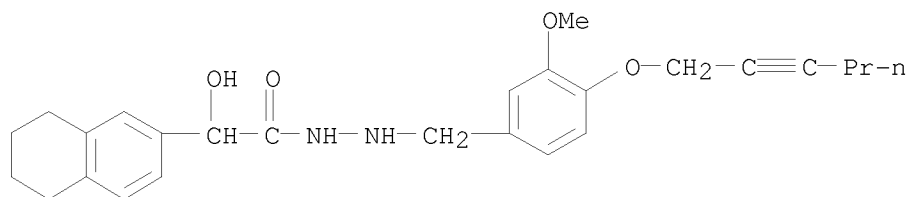
RN 1055765-40-6 CAPLUS

CN 2-Naphthaleneacetamide, N-[[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055765-41-7 CAPLUS

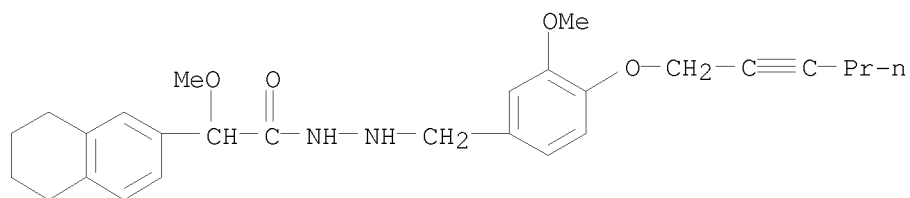
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-, 2-[[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055765-42-8 CAPLUS

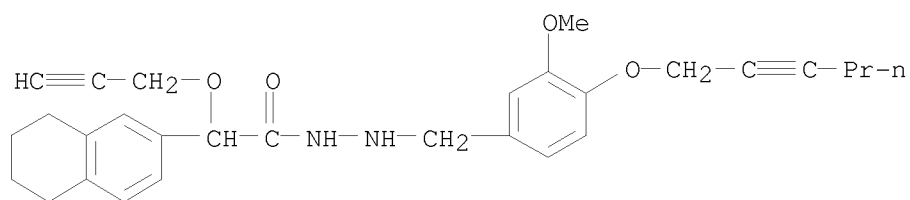
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CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-,
2-[[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)



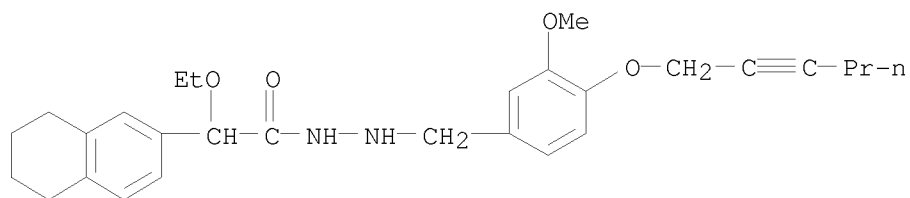
RN 1055765-43-9 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)



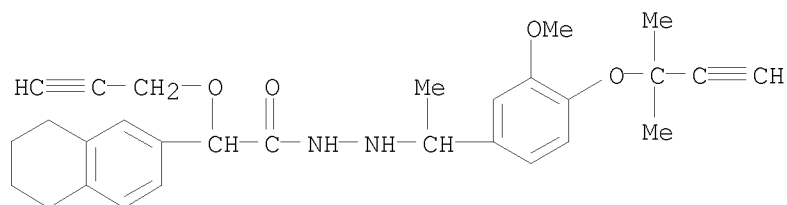
RN 1055765-44-0 CAPLUS

CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
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RN 1055766-37-4 CAPLUS

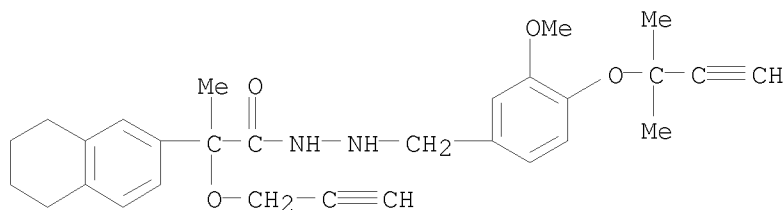
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[1-[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]hydrazide
(CA INDEX NAME)



RN 1055766-38-5 CAPLUS

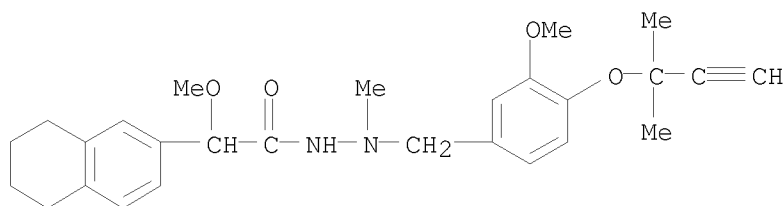
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CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)



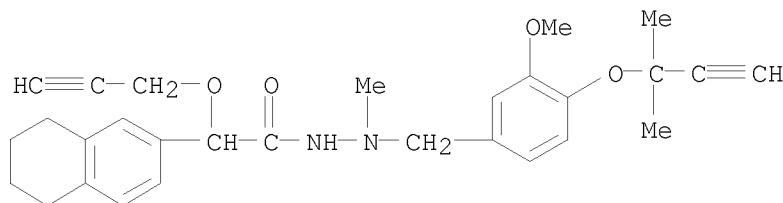
RN 1055766-39-6 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-, 2-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



RN 1055766-40-9 CAPLUS

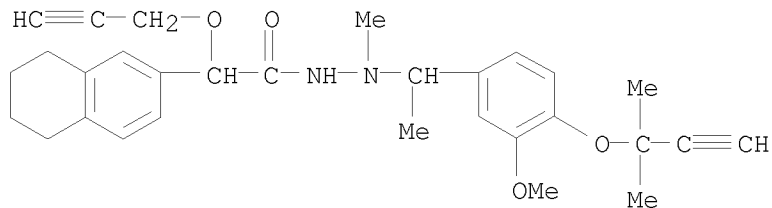
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RN 1055766-41-0 CAPLUS

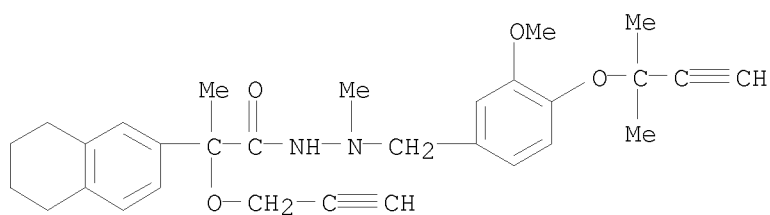
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10/513699



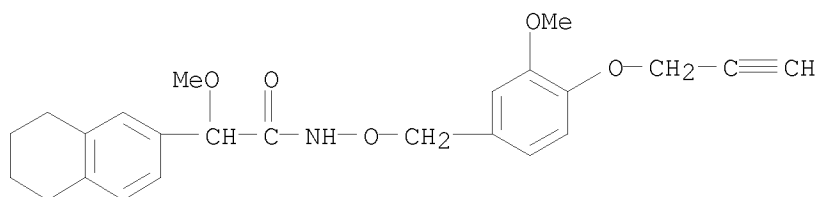
RN 1055766-42-1 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



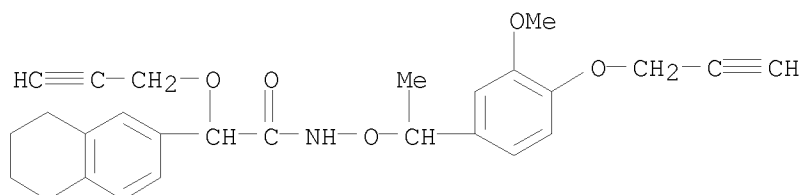
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CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -methoxy-N-[[3-methoxy-4-(2-propyn-1-yloxy)phenyl]methoxy]- (CA INDEX NAME)



RN 1055767-04-8 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[1-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethoxy]- α -(2-propyn-1-yloxy)- (CA INDEX NAME)

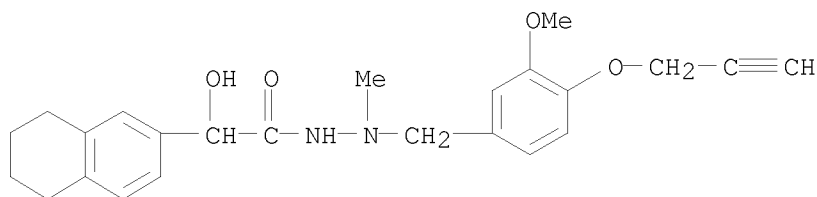


RN 1055767-05-9 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-, 2-[[3-methoxy-4-(2-propyn-1-yloxy)phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)

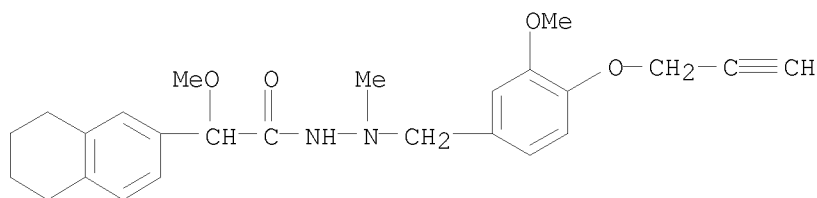
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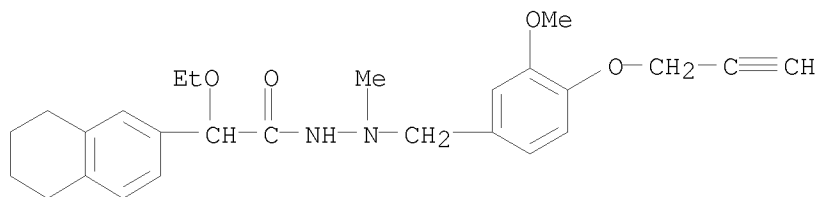
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CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-,
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INDEX NAME)



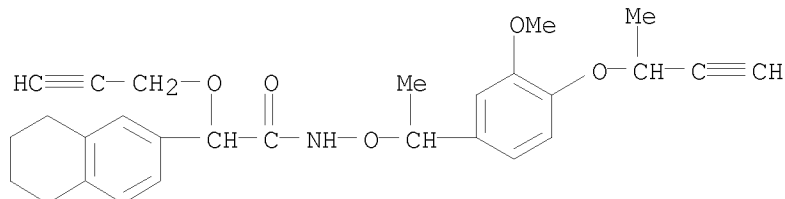
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CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
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INDEX NAME)



RN 1055767-31-1 CAPLUS

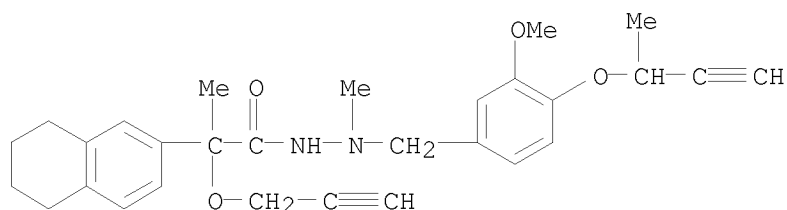
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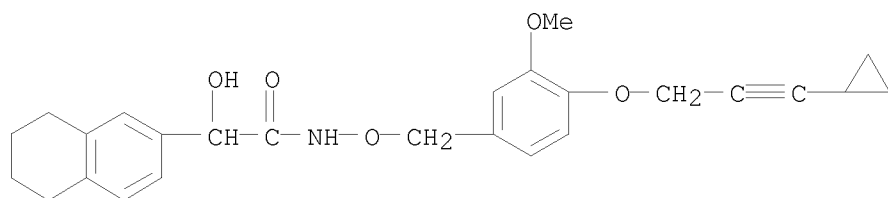
RN 1055767-32-2 CAPLUS

2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



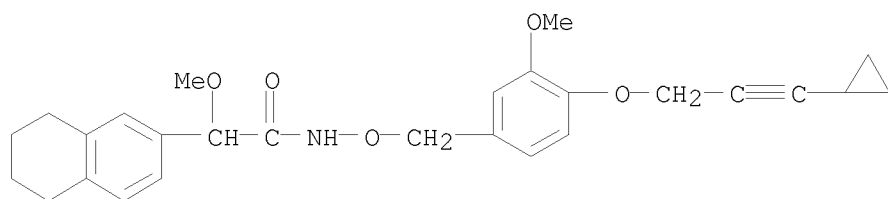
RN 1055768-24-5 CAPLUS

2-Naphthaleneacetamide, N-[[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro- α -hydroxy- (CA INDEX NAME)



RN 1055768-25-6 CAPLUS

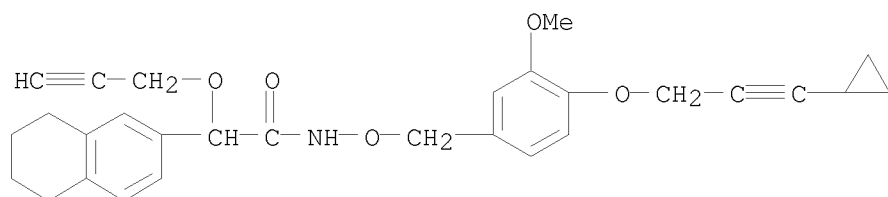
CN 2-Naphthaleneacetamide, N-[[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro- α -methoxy- (CA INDEX NAME)



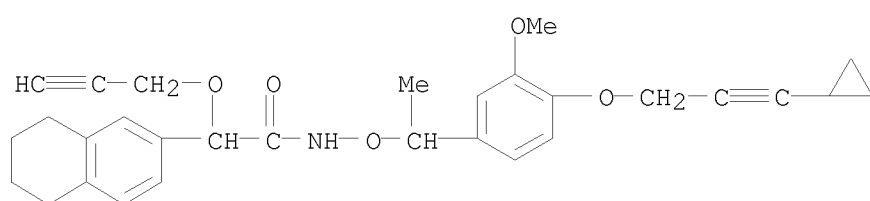
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CN 2-Naphthaleneacetamide, N-[[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)

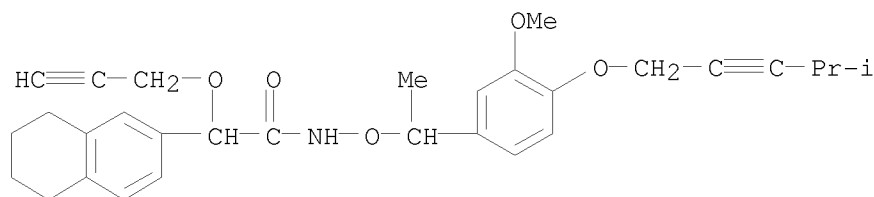
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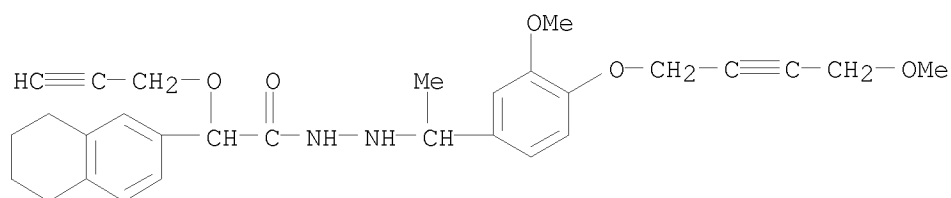
RN 1055768-27-8 CAPLUS
CN 2-Naphthaleneacetamide, N-[1-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethoxy]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055768-73-4 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

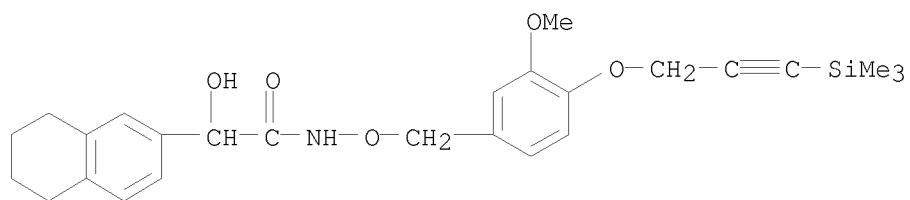


RN 1055769-27-1 CAPLUS
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[1-[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]ethyl]hydrazide (CA INDEX NAME)



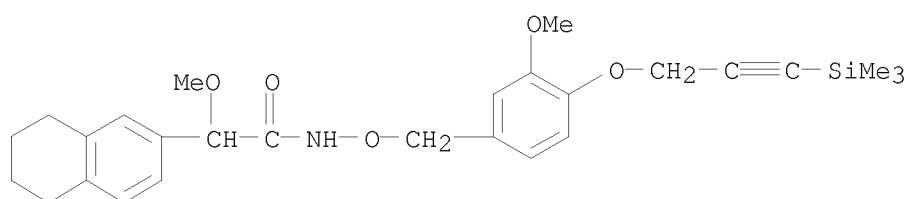
RN 1055769-95-3 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -hydroxy-N-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]methoxy]- (CA INDEX NAME)

10/513699



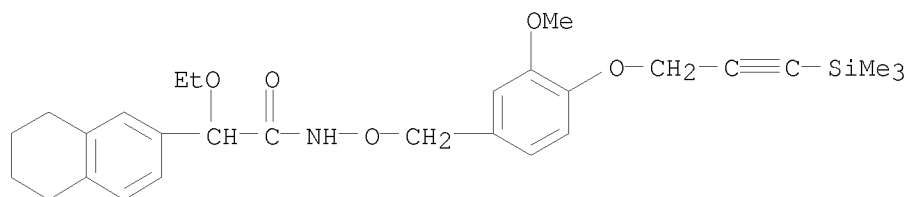
RN 1055769-96-4 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -methoxy-N-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]methoxy]- (CA INDEX NAME)



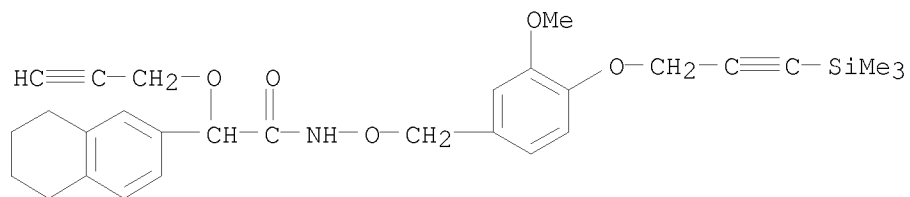
RN 1055769-97-5 CAPLUS

CN 2-Naphthaleneacetamide, α -ethoxy-5,6,7,8-tetrahydro-N-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]methoxy]- (CA INDEX NAME)



RN 1055769-98-6 CAPLUS

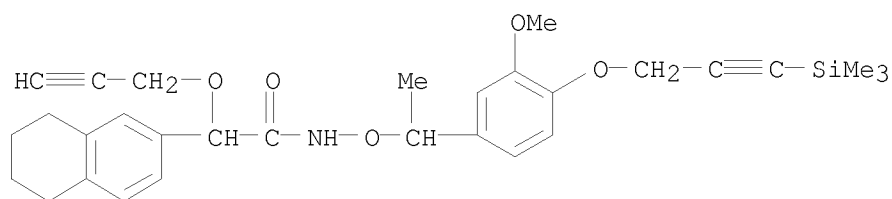
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]methoxy]- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055769-99-7 CAPLUS

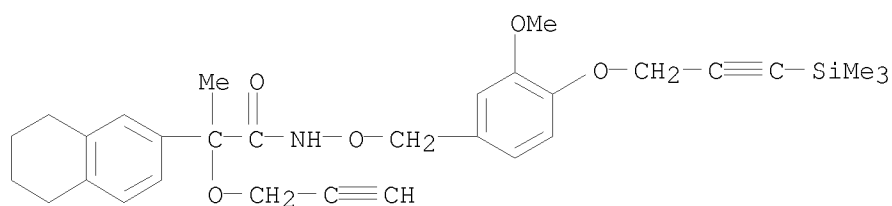
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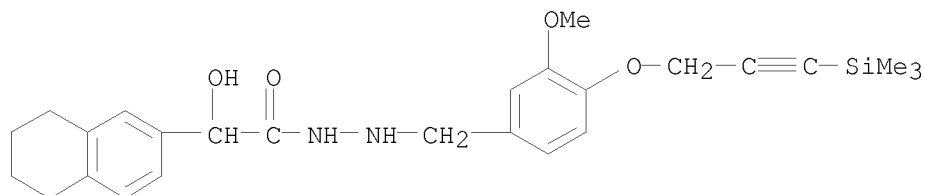
RN 1055770-00-7 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]methoxy]-α-methyl-α-(2-propyn-1-yloxy)- (CA INDEX NAME)



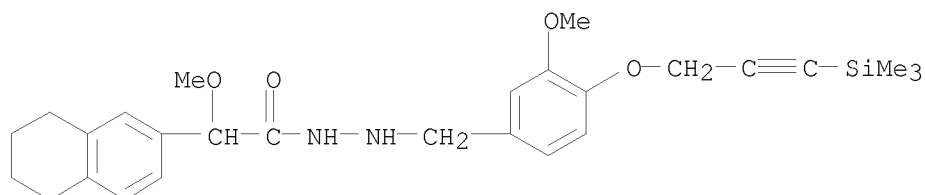
RN 1055770-01-8 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro-α-hydroxy-, 2-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055770-02-9 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro-α-methoxy-, 2-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]methyl]hydrazide (CA INDEX NAME)

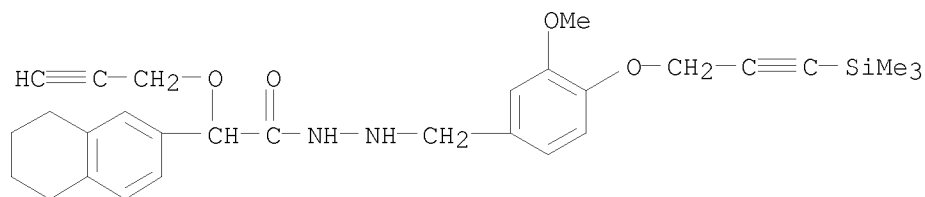


RN 1055770-03-0 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro-α-(2-propyn-1-yloxy)-, 2-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]methyl]hydrazide (CA INDEX NAME)

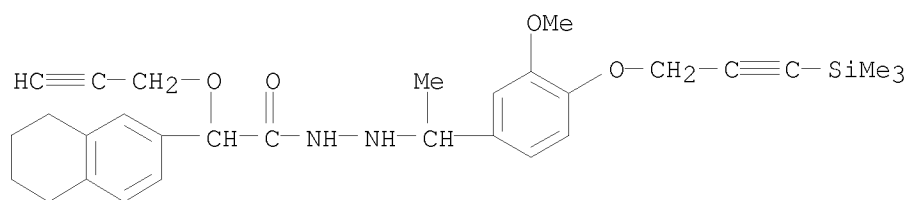
10/513699

yl]oxy]phenyl]methyl]hydrazide (CA INDEX NAME)



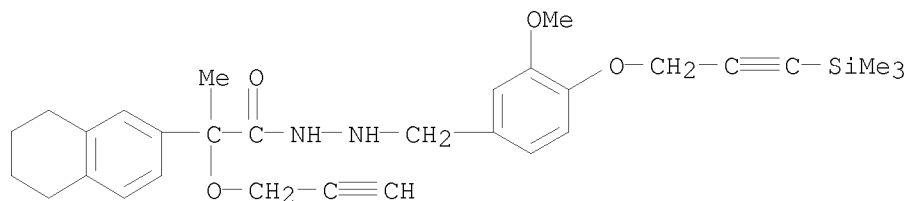
RN 1055770-04-1 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[1-[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]ethyl]hydrazide (CA INDEX NAME)



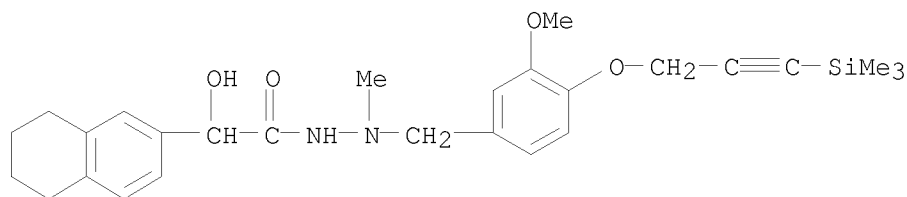
RN 1055770-05-2 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055770-06-3 CAPLUS

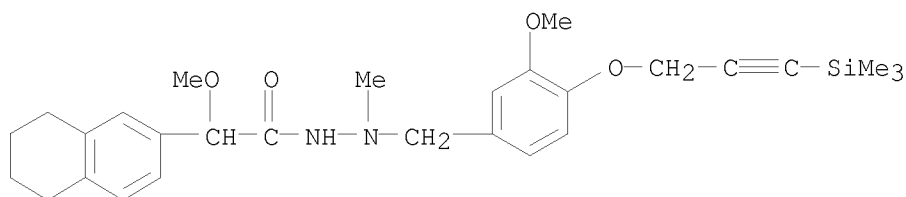
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-, 2-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



10/513699

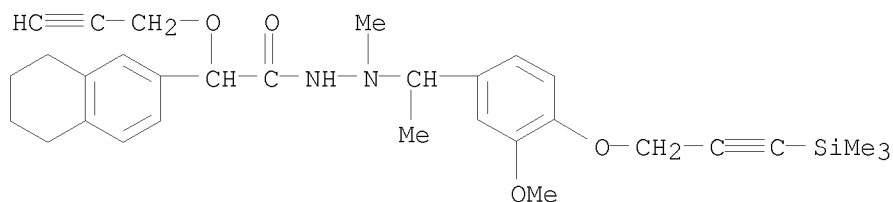
RN 1055770-07-4 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-,
2-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]methyl]-2-
methylhydrazide (CA INDEX NAME)



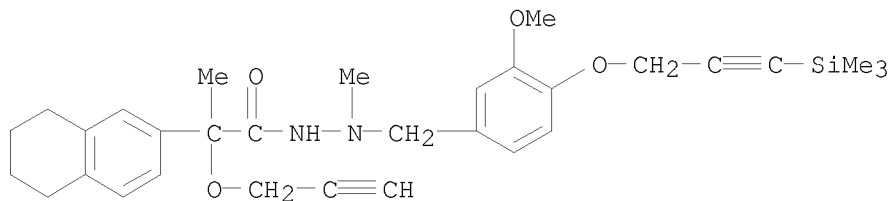
RN 1055770-08-5 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
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methylhydrazide (CA INDEX NAME)



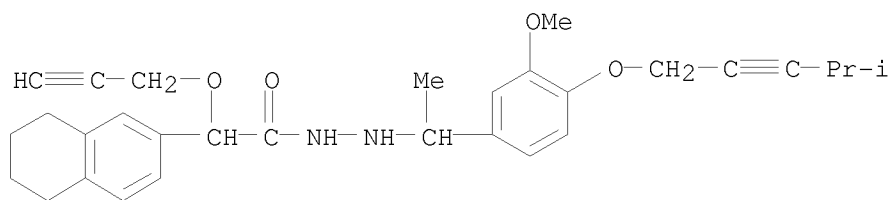
RN 1055770-09-6 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-
propyn-1-yloxy)-, 2-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-
yl]oxy]phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



RN 1055770-97-2 CAPLUS

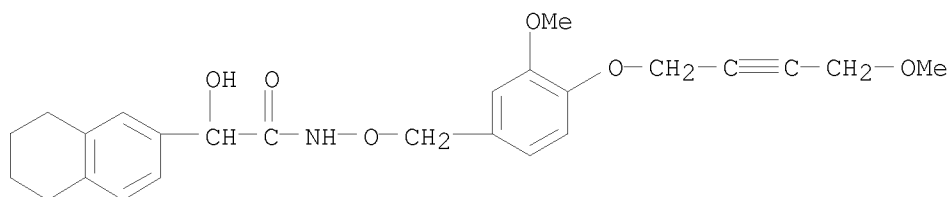
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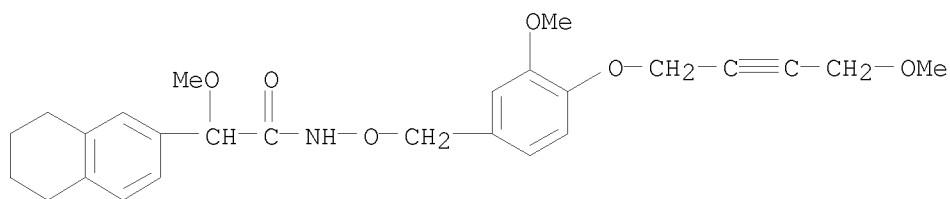
RN 1055771-22-6 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -hydroxy-N-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methoxy]- (CA INDEX NAME)



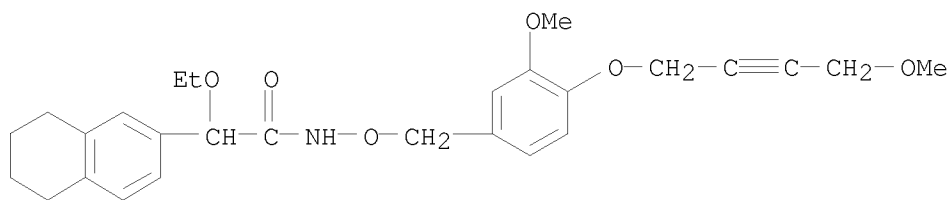
RN 1055771-23-7 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -methoxy-N-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methoxy]- (CA INDEX NAME)



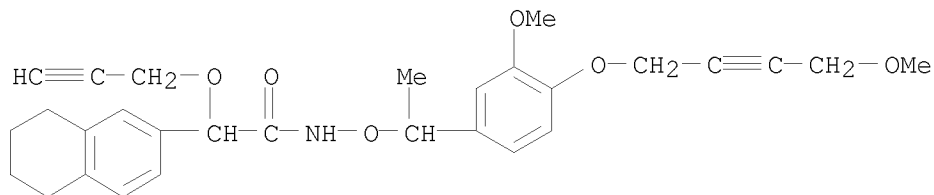
RN 1055771-24-8 CAPLUS

CN 2-Naphthaleneacetamide, α -ethoxy-5,6,7,8-tetrahydro-N-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methoxy]- (CA INDEX NAME)



RN 1055771-25-9 CAPLUS

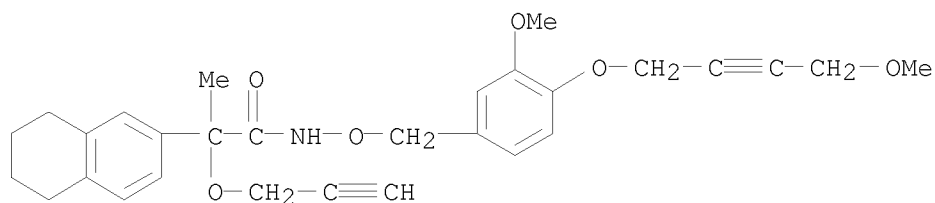
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[1-[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]ethoxy]- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



10/513699

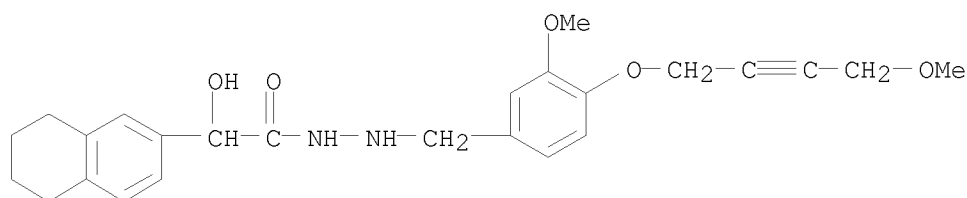
RN 1055771-26-0 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methoxy]- α -methyl- α -(2-propyn-1-yloxy)-
(CA INDEX NAME)



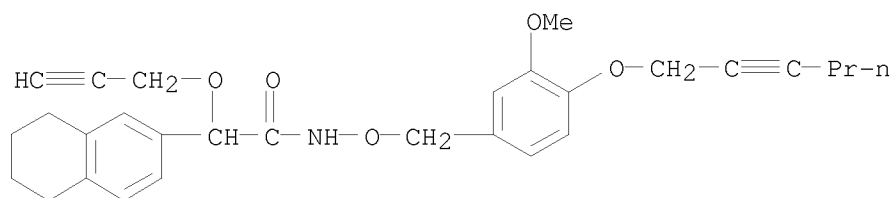
RN 1055771-27-1 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-,
2-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methyl]hydrazide (CA
INDEX NAME)



RN 1055771-91-9 CAPLUS

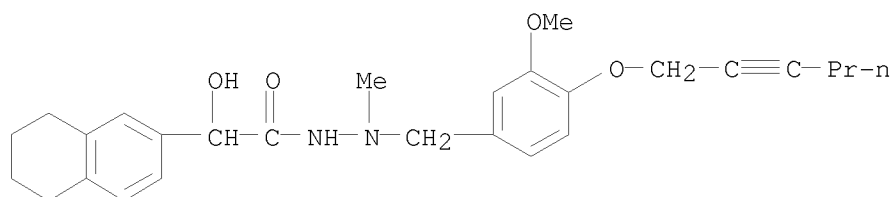
CN 2-Naphthaleneacetamide, N-[[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]methoxy]-
5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055772-00-3 CAPLUS

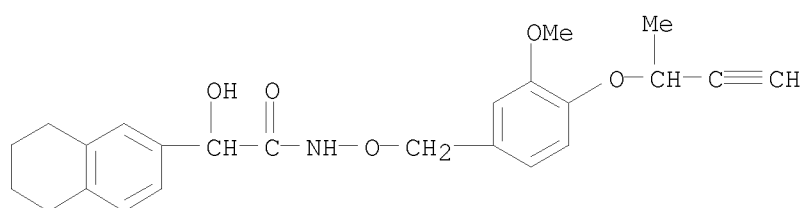
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-,
2-[[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]methyl]-2-methylhydrazide (CA
INDEX NAME)

10/513699



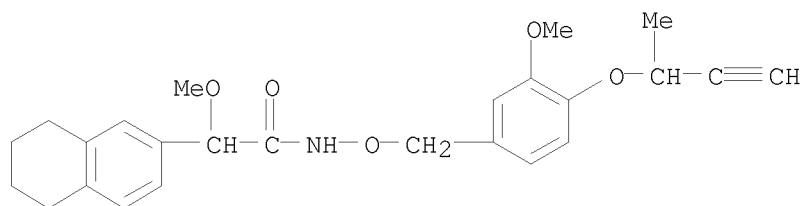
RN 1055772-94-5 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -hydroxy-N-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methoxy]- (CA INDEX NAME)



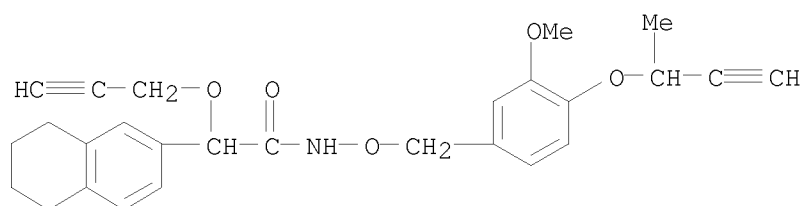
RN 1055772-95-6 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -methoxy-N-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methoxy]- (CA INDEX NAME)



RN 1055772-96-7 CAPLUS

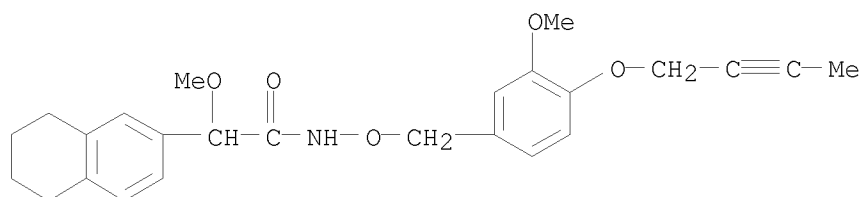
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methoxy]- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055773-09-5 CAPLUS

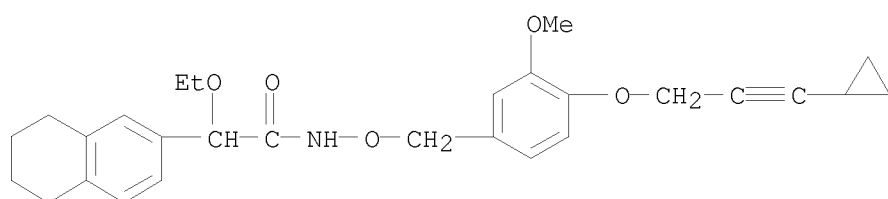
CN 2-Naphthaleneacetamide, N-[[4-(2-butyn-1-yloxy)-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro- α -methoxy- (CA INDEX NAME)

10/513699



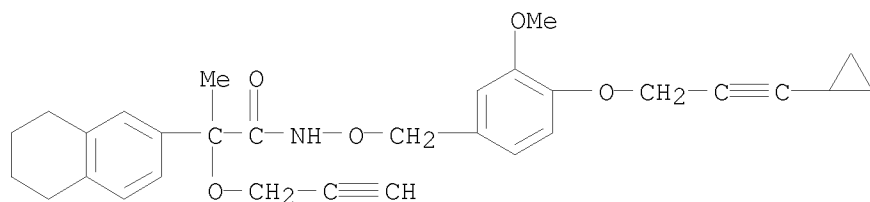
RN 1055774-87-2 CAPLUS

CN 2-Naphthaleneacetamide, N-[[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methoxy]-α-ethoxy-5,6,7,8-tetrahydro- (CA INDEX NAME)



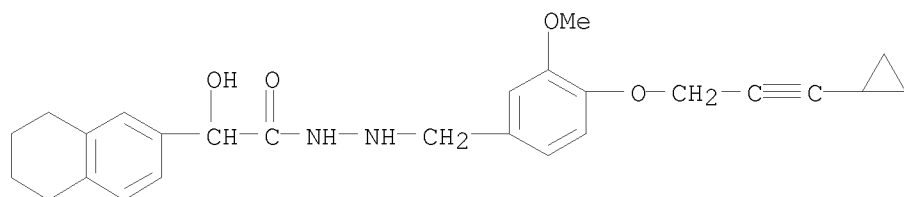
RN 1055774-89-4 CAPLUS

CN 2-Naphthaleneacetamide, N-[[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro-α-methyl-α-(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055774-90-7 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro-α-hydroxy-, 2-[[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)

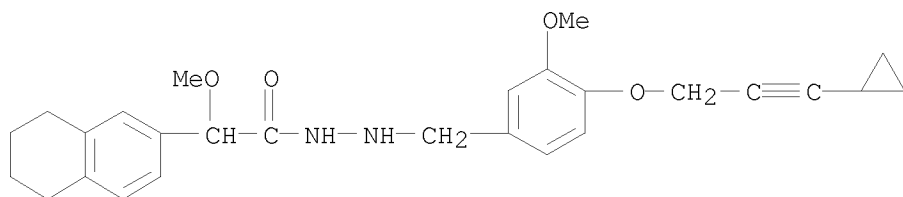


RN 1055774-91-8 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro-α-methoxy-, 2-[[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]hydrazide

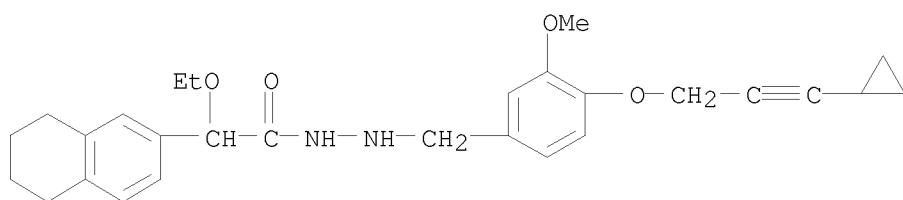
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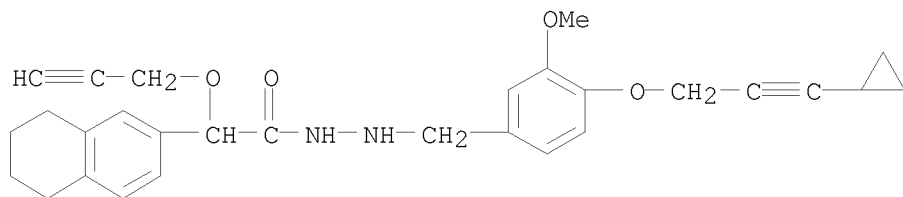
RN 1055774-92-9 CAPLUS

CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
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(CA INDEX NAME)



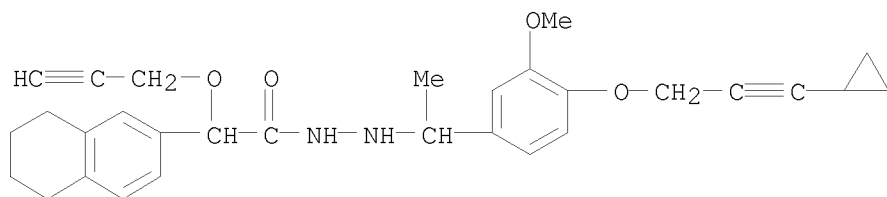
RN 1055774-93-0 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
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(CA INDEX NAME)



RN 1055774-94-1 CAPLUS

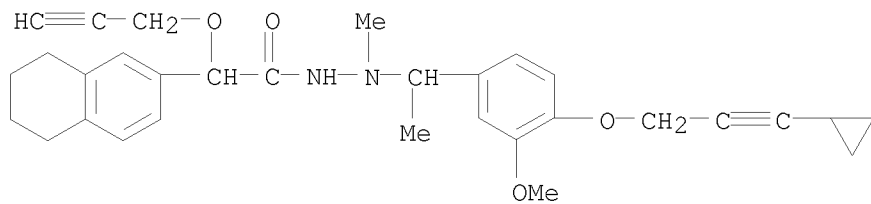
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
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(CA INDEX NAME)



10/513699

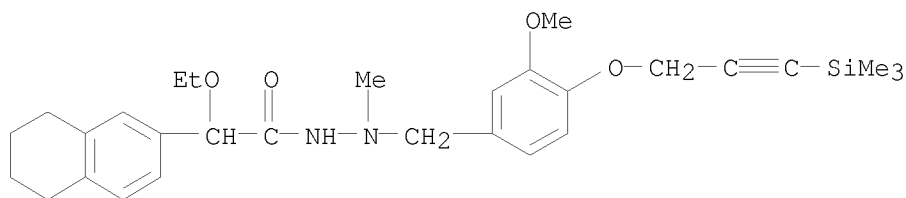
RN 1055774-95-2 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
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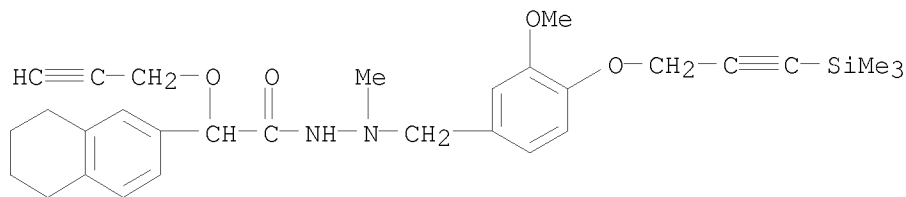
RN 1055775-30-8 CAPLUS

CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
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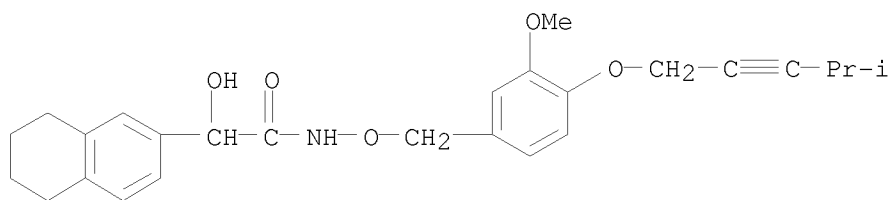
RN 1055775-31-9 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
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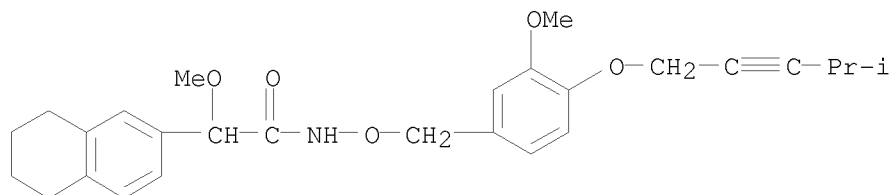
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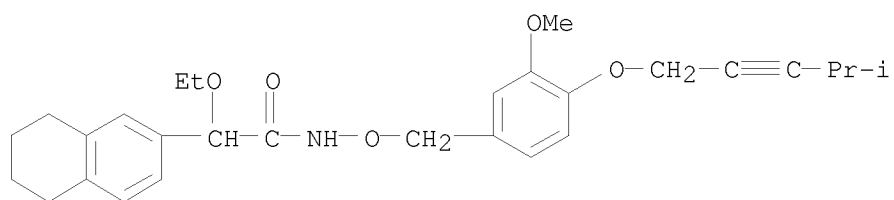


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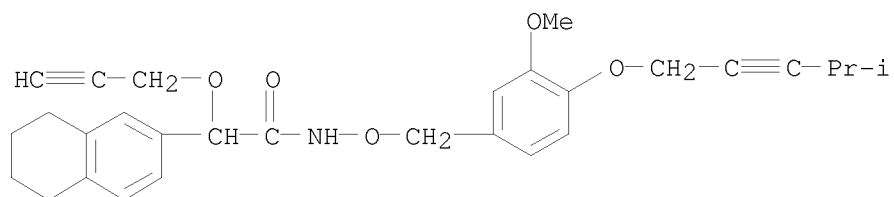
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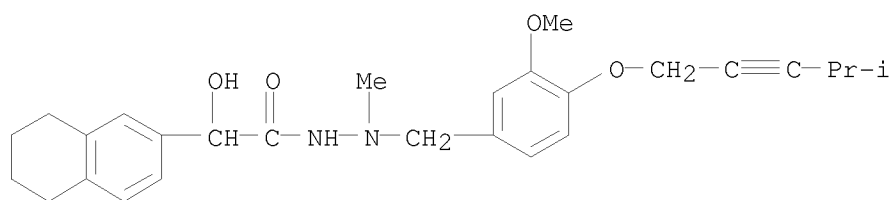
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RN 1055775-72-8 CAPLUS
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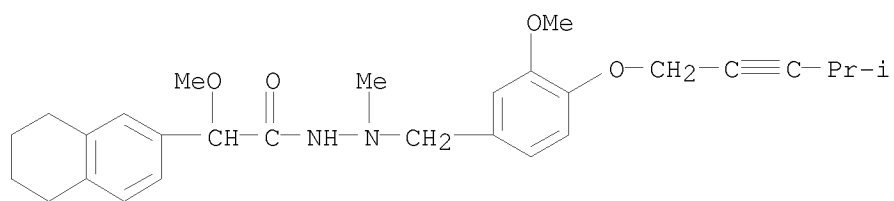


RN 1055775-73-9 CAPLUS
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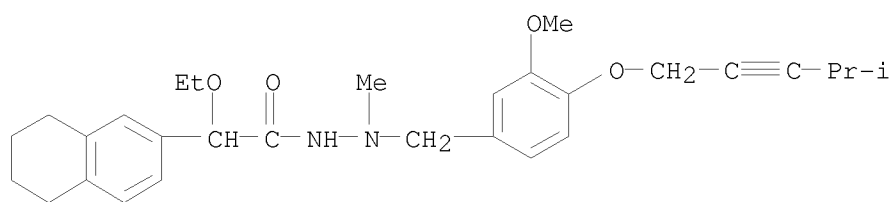


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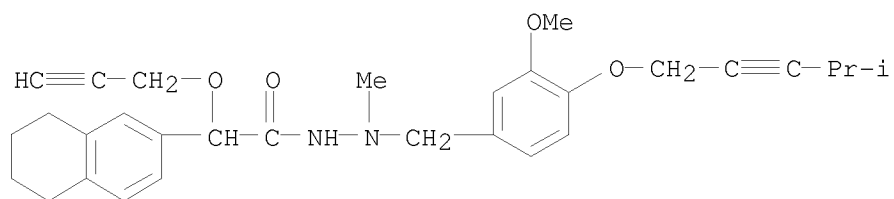
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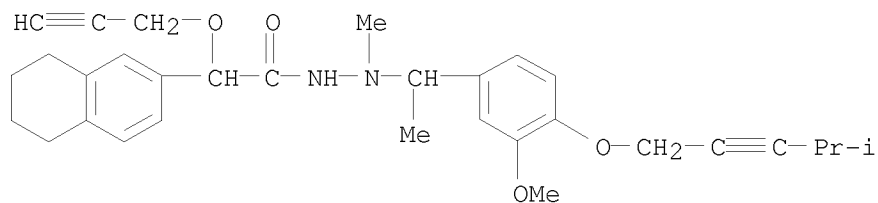
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RN 1055775-76-2 CAPLUS
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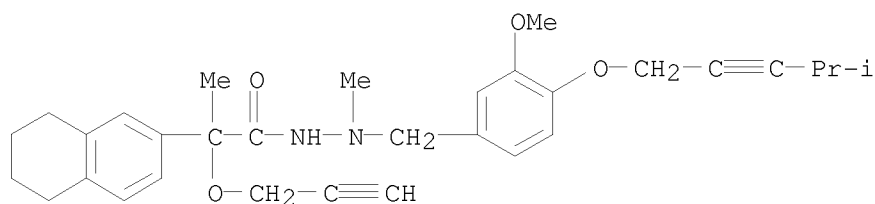


RN 1055775-77-3 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



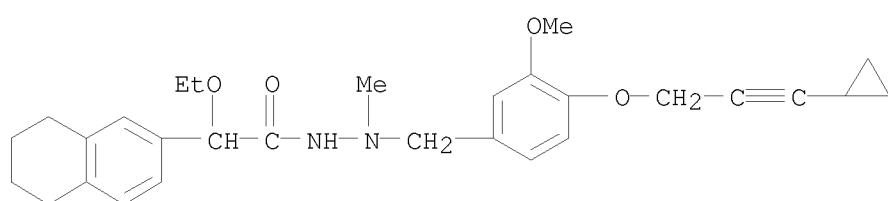
RN 1055775-78-4 CAPLUS
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10/513699



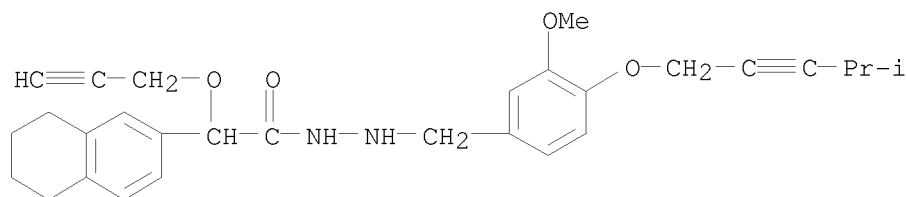
RN 1055775-89-7 CAPLUS

CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
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methylhydrazide (CA INDEX NAME)



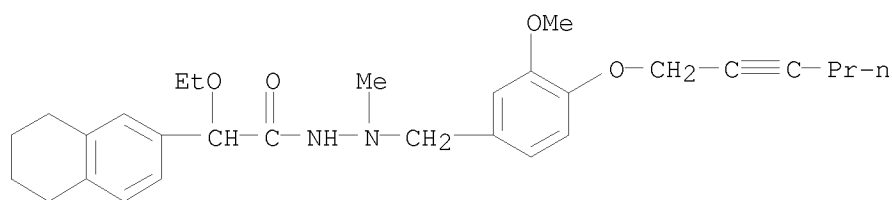
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CN INDEX NAME NOT YET ASSIGNED



RN 1055776-74-3 CAPLUS

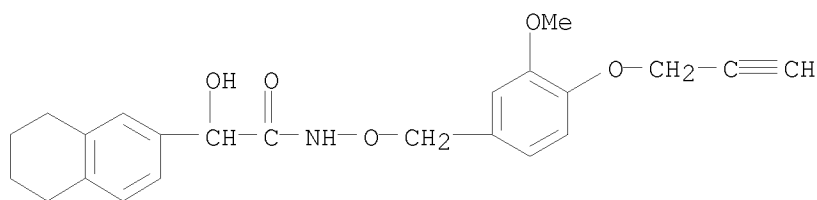
CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
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INDEX NAME)



RN 1055776-95-8 CAPLUS

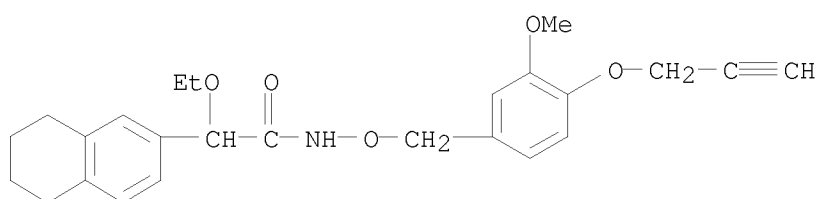
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -hydroxy-N-[[3-methoxy-4-
(2-propyn-1-yloxy)phenyl]methoxy]- (CA INDEX NAME)

10/513699



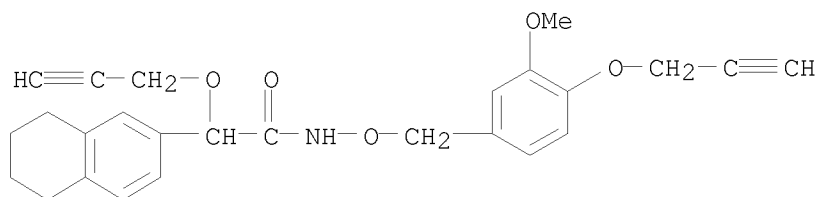
RN 1055776-96-9 CAPLUS

CN 2-Naphthaleneacetamide, α -ethoxy-5,6,7,8-tetrahydro-N-[[3-methoxy-4-(2-propyn-1-yloxy)phenyl]methoxy]- (CA INDEX NAME)



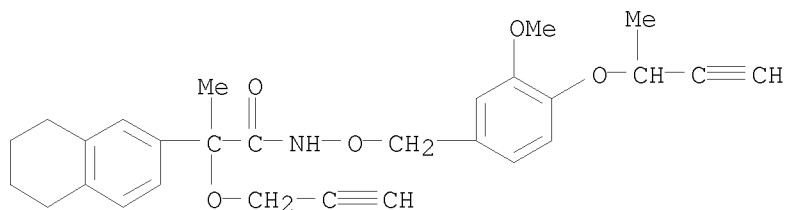
RN 1055776-97-0 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[[3-methoxy-4-(2-propyn-1-yloxy)phenyl]methoxy]- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055777-87-1 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methoxy]- α -methyl- α -(2-propyn-1-yloxy)- (CA INDEX NAME)

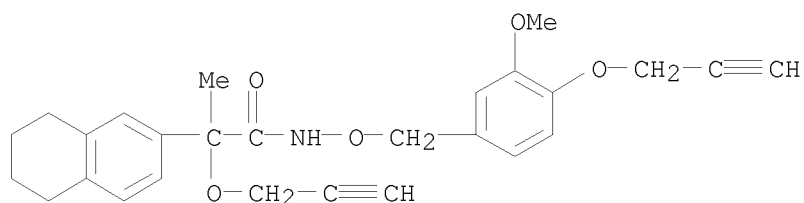


RN 1055777-95-1 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[[3-methoxy-4-(2-propyn-1-yloxy)phenyl]methoxy]- α -methyl- α -(2-propyn-1-yloxy)- (CA

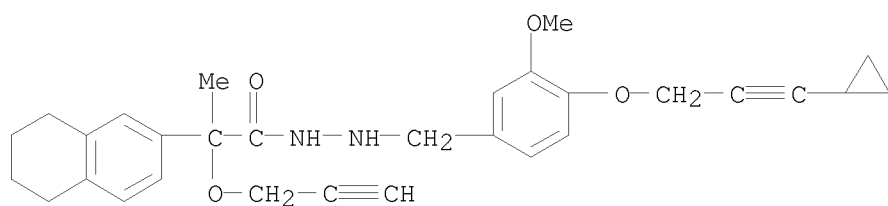
10/513699

INDEX NAME)



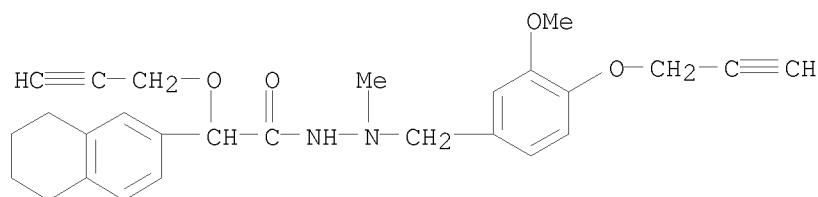
RN 1055778-06-7 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)



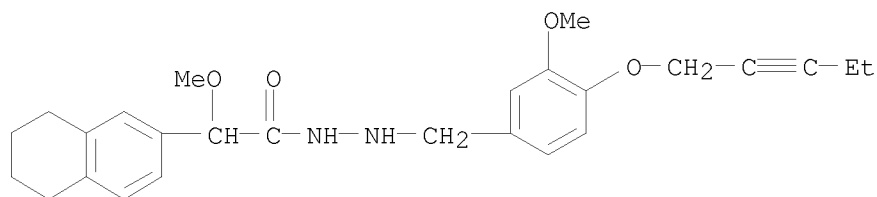
RN 1055778-37-4 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[[3-methoxy-4-(2-propyn-1-yloxy)phenyl]methyl]-2-methylhydrazide (CA
INDEX NAME)



RN 1055778-58-9 CAPLUS

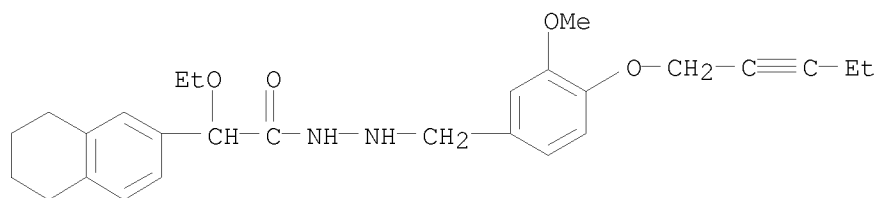
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-,
2-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055778-59-0 CAPLUS

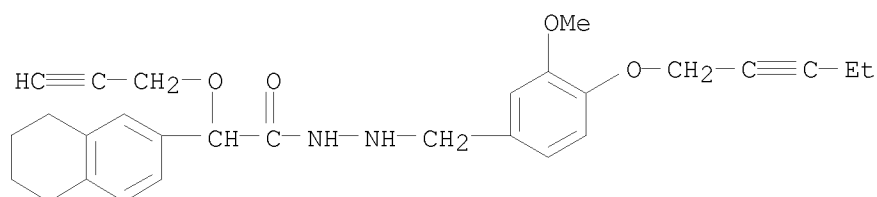
10/513699

CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
2-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methyl]hydrazide (CA INDEX NAME)



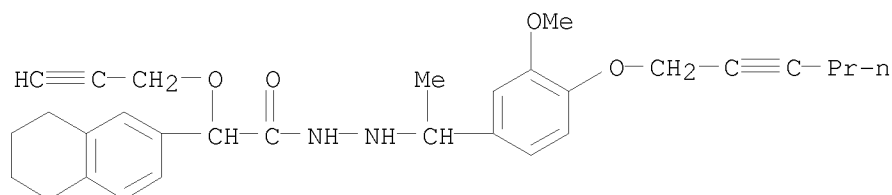
RN 1055778-60-3 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methyl]hydrazide (CA INDEX NAME)



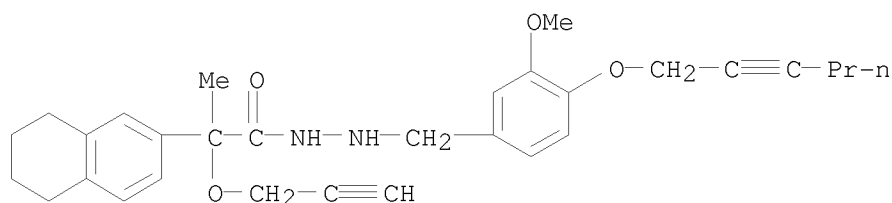
RN 1055778-70-5 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[1-[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]ethyl]hydrazide (CA INDEX NAME)



RN 1055778-71-6 CAPLUS

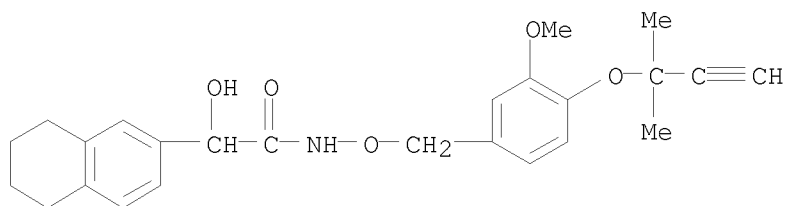
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-,
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RN 1055779-76-4 CAPLUS

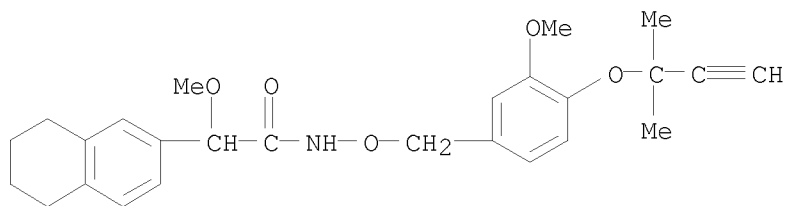
10/513699

CN 2-Naphthaleneacetamide, N-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro- α -hydroxy- (CA INDEX NAME)



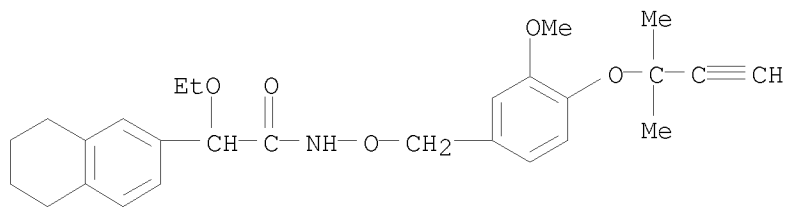
RN 1055779-77-5 CAPLUS

CN 2-Naphthaleneacetamide, N-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro- α -methoxy- (CA INDEX NAME)



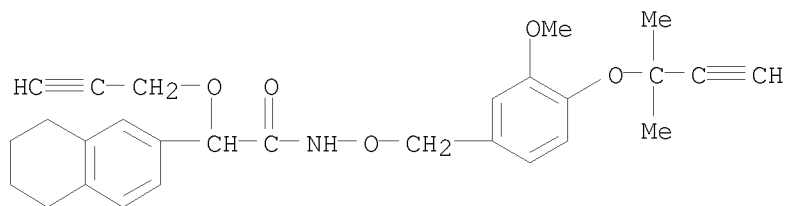
RN 1055779-78-6 CAPLUS

CN 2-Naphthaleneacetamide, N-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methoxy]- α -ethoxy-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 1055779-79-7 CAPLUS

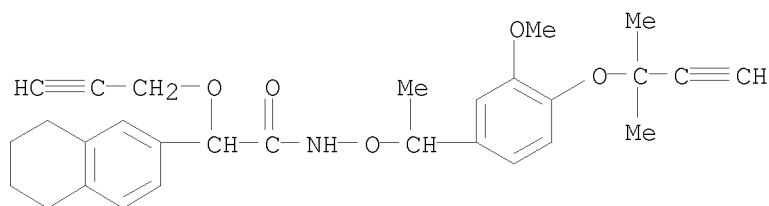
CN 2-Naphthaleneacetamide, N-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



10/513699

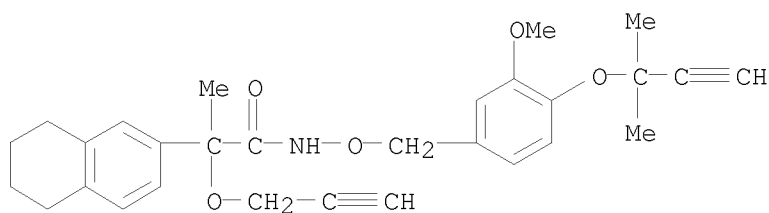
RN 1055779-80-0 CAPLUS

CN 2-Naphthaleneacetamide, N-[1-[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethoxy]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



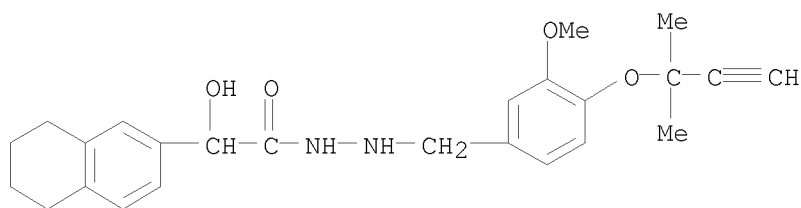
RN 1055779-81-1 CAPLUS

CN 2-Naphthaleneacetamide, N-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055779-82-2 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-, 2-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)



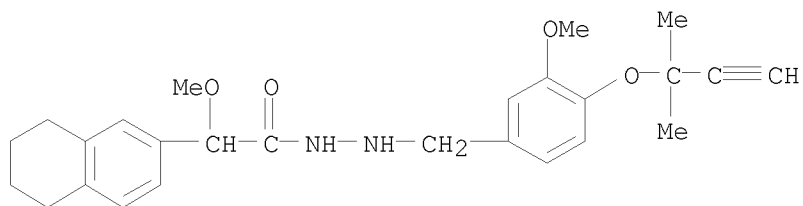
RN 1055779-83-3 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-, 2-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)

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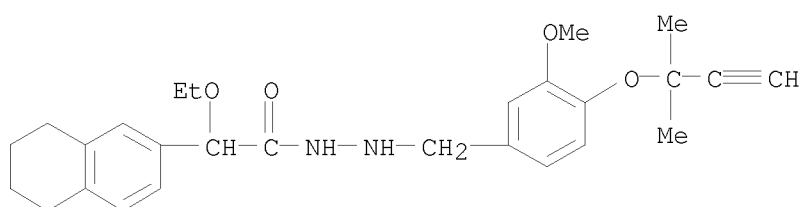
Erich Leese

10/513699



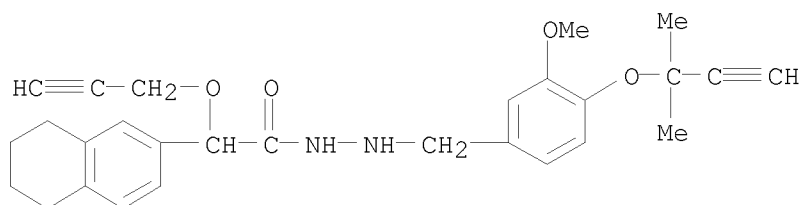
RN 1055779-84-4 CAPLUS

CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
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(CA INDEX NAME)



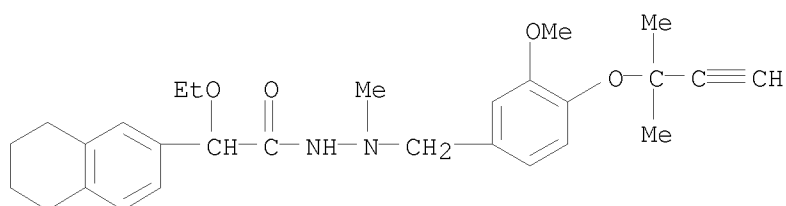
RN 1055779-85-5 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]hydrazide
(CA INDEX NAME)



RN 1055779-86-6 CAPLUS

CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
2-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]-2-
methylhydrazide (CA INDEX NAME)

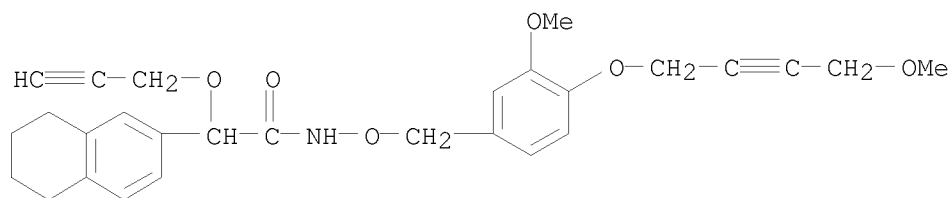


RN 1055780-35-2 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[[3-methoxy-4-[(4-methoxy-2-
ethoxy-1-propyn-1-yl)oxy]phenyl]methyl] (CA INDEX NAME)

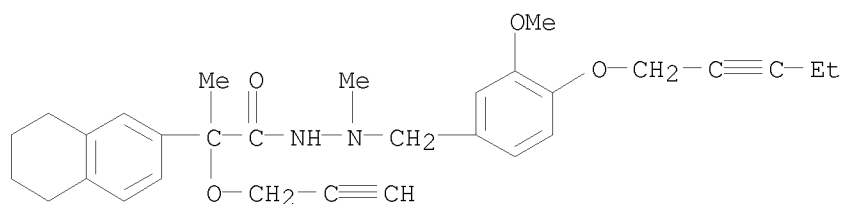
10/513699

butyn-1-yl)oxy]phenyl]methoxy]- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



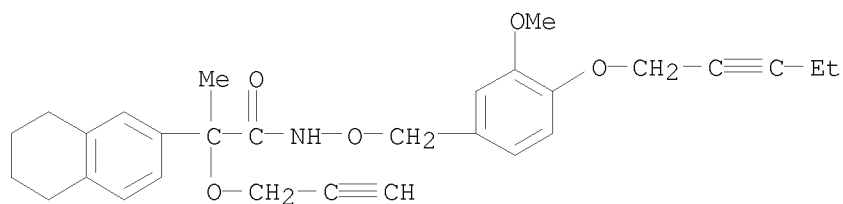
RN 1055780-93-2 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



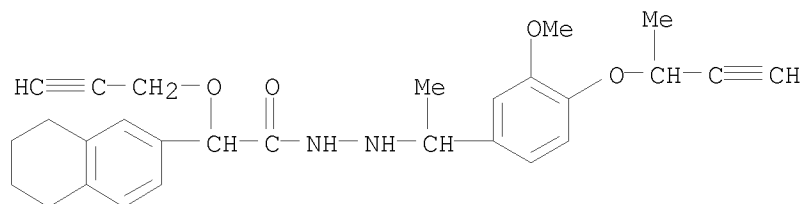
RN 1055781-13-9 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methoxy]- α -methyl- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055782-11-0 CAPLUS

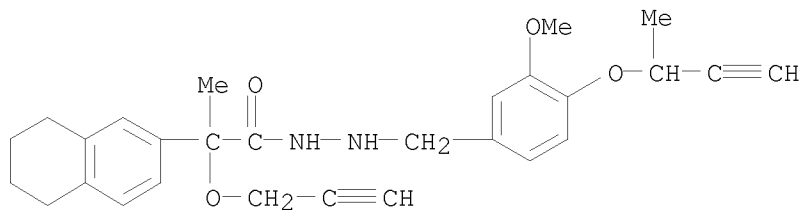
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[1-[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]ethyl]hydrazide (CA INDEX NAME)



10/513699

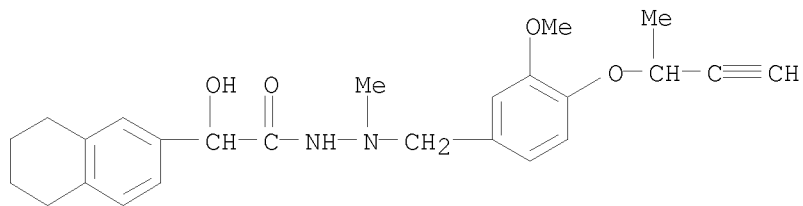
RN 1055782-12-1 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methyl]hydrazide (CA INDEX NAME)



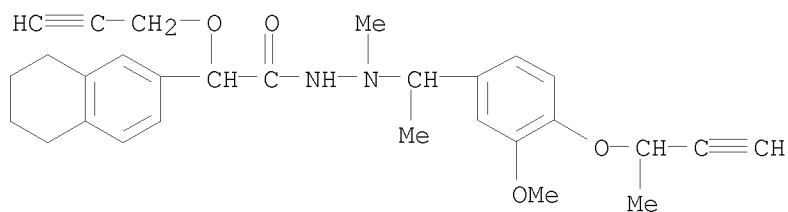
RN 1055782-13-2 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-, 2-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



RN 1055782-14-3 CAPLUS

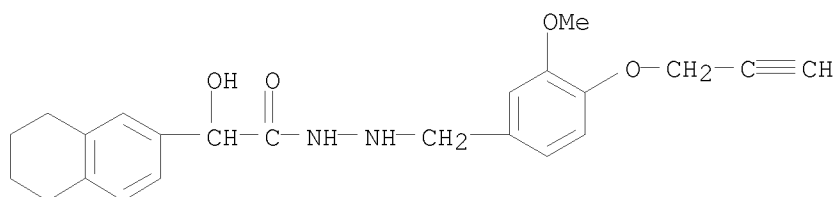
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[1-[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]ethyl]-2-methylhydrazide (CA INDEX NAME)



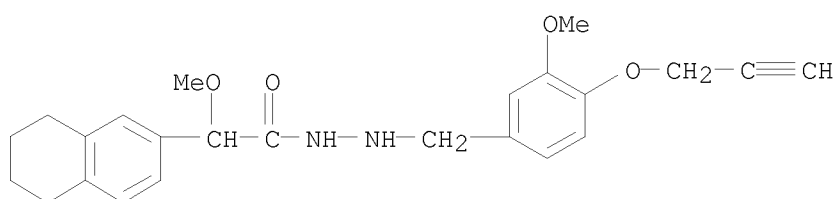
RN 1055782-36-9 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-, 2-[[3-methoxy-4-(2-propyn-1-yloxy)phenyl]methyl]hydrazide (CA INDEX NAME)

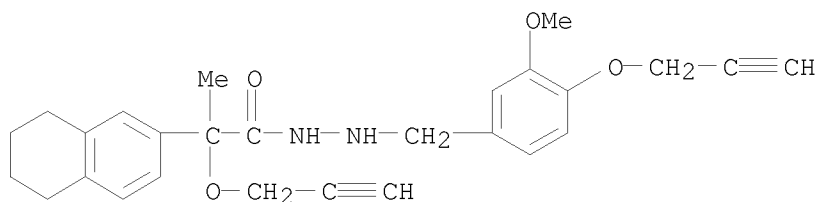
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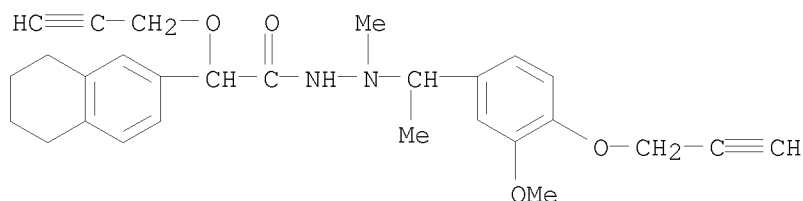
RN 1055782-37-0 CAPLUS
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-,
2-[[3-methoxy-4-(2-propyn-1-yloxy)phenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055782-38-1 CAPLUS
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[3-methoxy-4-(2-propyn-1-yloxy)phenyl]methyl]hydrazide (CA INDEX NAME)



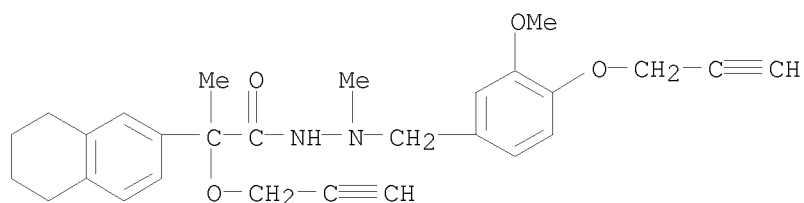
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CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[1-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-2-methylhydrazide (CA INDEX NAME)



RN 1055782-40-5 CAPLUS
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[3-methoxy-4-(2-propyn-1-yloxy)phenyl]methyl]-2-

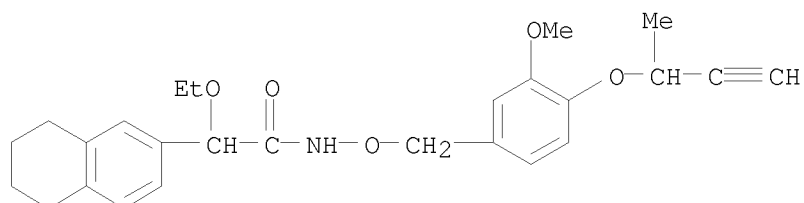
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methylhydrazide (CA INDEX NAME)



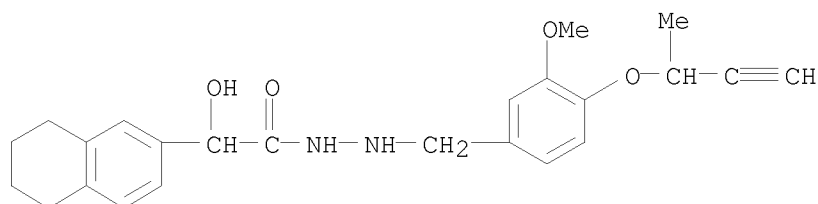
RN 1055782-91-6 CAPLUS

CN 2-Naphthaleneacetamide, α -ethoxy-5,6,7,8-tetrahydro-N-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methoxy]- (CA INDEX NAME)



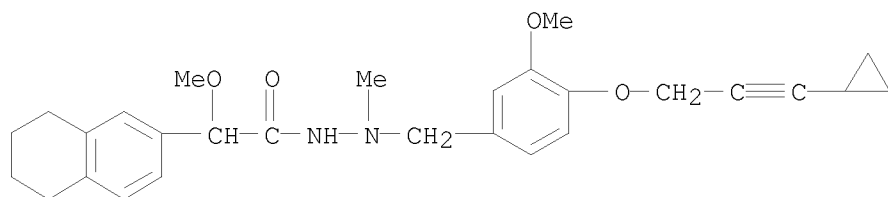
RN 1055782-92-7 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-, 2-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055783-85-1 CAPLUS

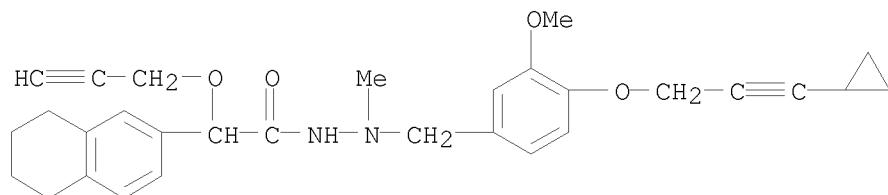
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-, 2-[[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



RN 1055783-86-2 CAPLUS

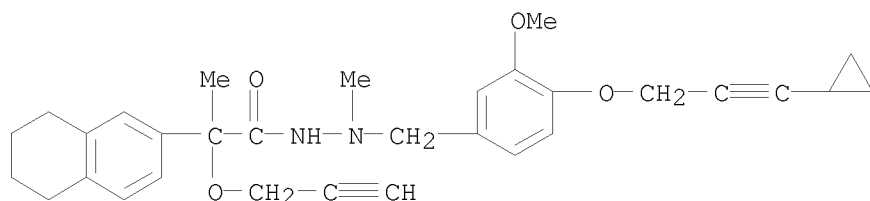
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CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



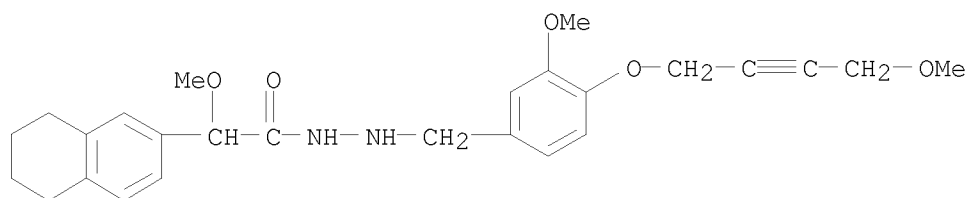
RN 1055783-87-3 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



RN 1055784-30-9 CAPLUS

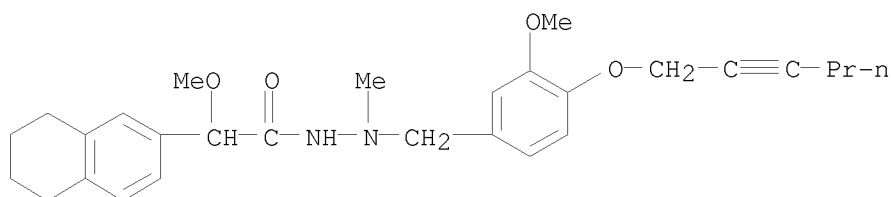
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-, 2-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055784-57-0 CAPLUS

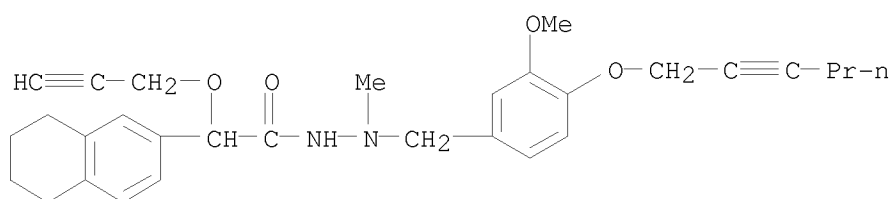
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-, 2-[[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]methyl]-2-methylhydrazide (CA INDEX NAME)

10/513699



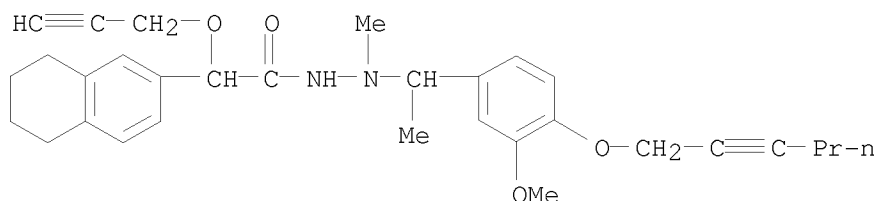
RN 1055784-58-1 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



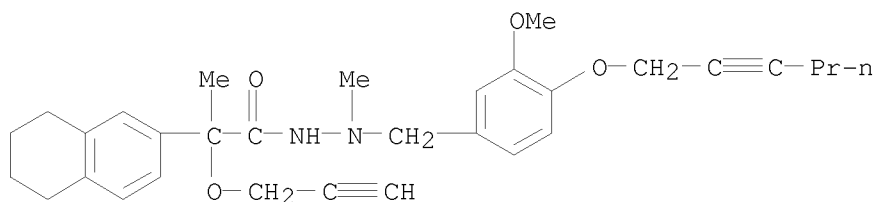
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CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[1-[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]ethyl]-2-methylhydrazide (CA INDEX NAME)



RN 1055784-60-5 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]methyl]-2-methylhydrazide (CA INDEX NAME)

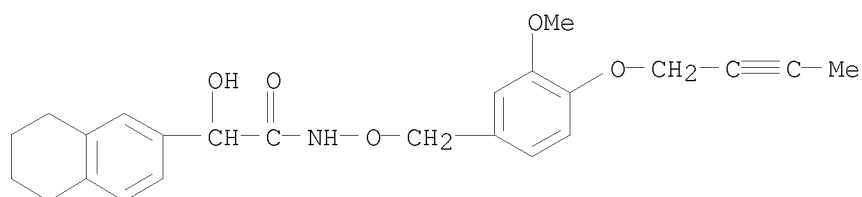


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CN 2-Naphthaleneacetamide, N-[[4-(2-butyn-1-yloxy)-3-methoxyphenyl]methoxy]-

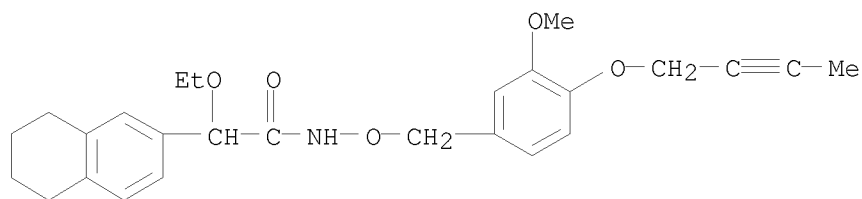
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5,6,7,8-tetrahydro- α -hydroxy- (CA INDEX NAME)



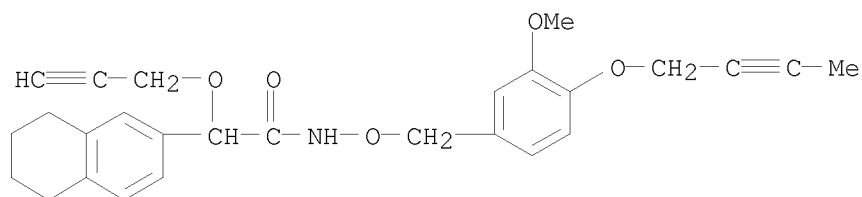
RN 1055784-90-1 CAPLUS

CN 2-Naphthaleneacetamide, N-[[4-(2-butyn-1-yloxy)-3-methoxyphenyl]methoxy]- α -ethoxy-5,6,7,8-tetrahydro- (CA INDEX NAME)



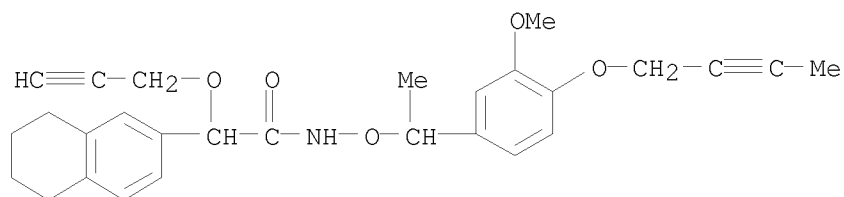
RN 1055784-91-2 CAPLUS

CN 2-Naphthaleneacetamide, N-[[4-(2-butyn-1-yloxy)-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055784-92-3 CAPLUS

CN 2-Naphthaleneacetamide, N-[1-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethoxy]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)

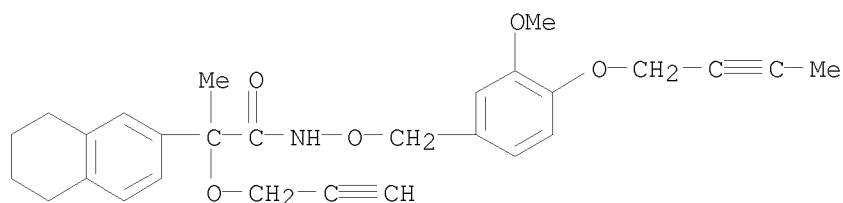


RN 1055784-93-4 CAPLUS

CN 2-Naphthaleneacetamide, N-[[4-(2-butyn-1-yloxy)-3-methoxyphenyl]methoxy]-5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)- (CA INDEX NAME)

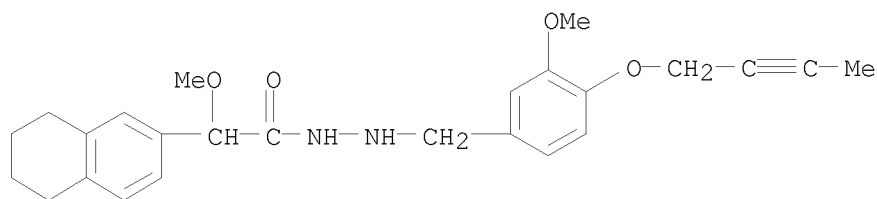
10/513699

NAME)



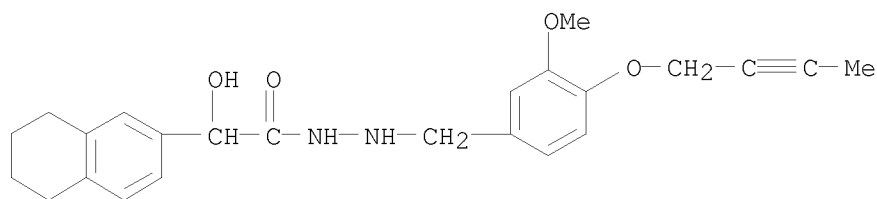
RN 1055784-94-5 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-,
2-[[4-(2-butyne-1-yloxy)-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)



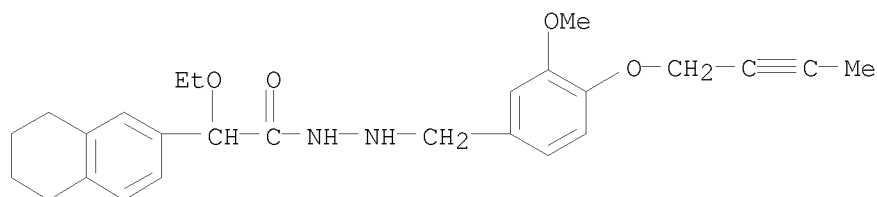
RN 1055784-95-6 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-,
2-[[4-(2-butyne-1-yloxy)-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055784-96-7 CAPLUS

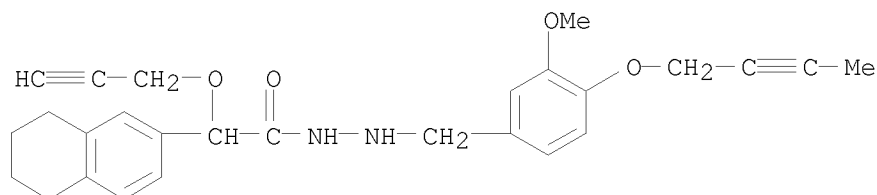
CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
2-[[4-(2-butyne-1-yloxy)-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055784-97-8 CAPLUS

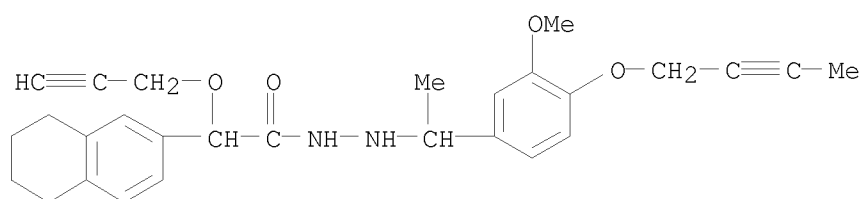
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[[4-(2-butyne-1-yloxy)-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)

10/513699



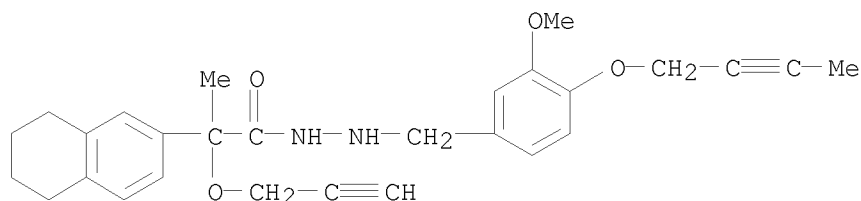
RN 1055784-98-9 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[1-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]hydrazide (CA INDEX NAME)



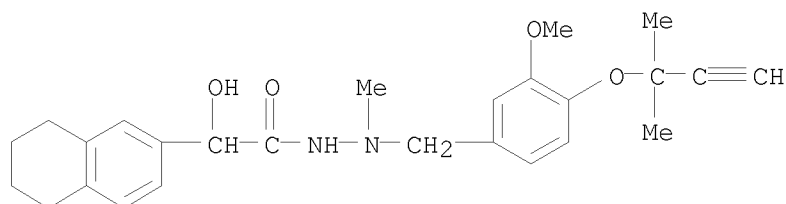
RN 1055784-99-0 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[4-(2-butyn-1-yloxy)-3-methoxyphenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055785-73-3 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-, 2-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]-2-methylhydrazide (CA INDEX NAME)

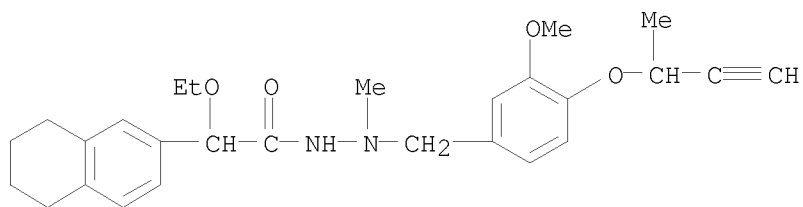


RN 1055786-75-8 CAPLUS

CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-, 2-[[4-[(1,1-dimethyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]-2-methylhydrazide (CA INDEX NAME)

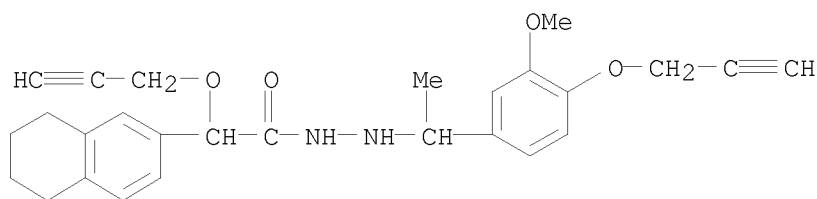
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2-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



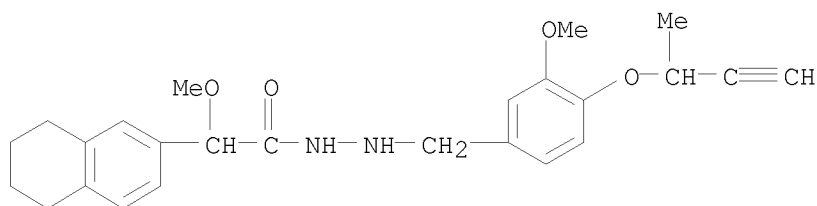
RN 1055787-00-2 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[1-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]hydrazide (CA INDEX NAME)



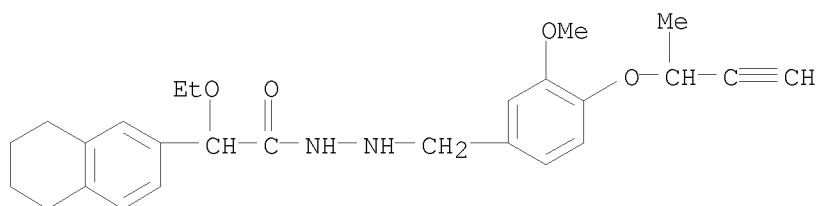
RN 1055787-60-4 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-, 2-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055787-61-5 CAPLUS

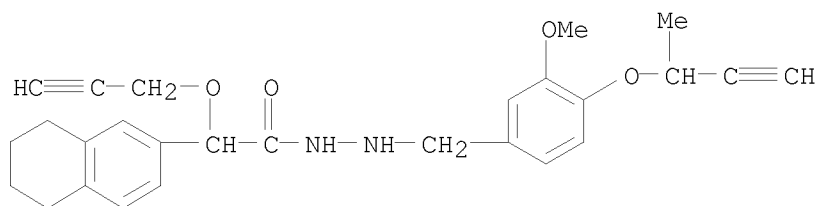
CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-, 2-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methyl]hydrazide (CA INDEX NAME)



10/513699

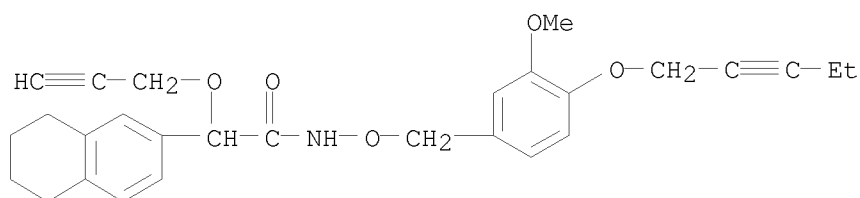
RN 1055787-62-6 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methyl]hydrazide (CA
INDEX NAME)



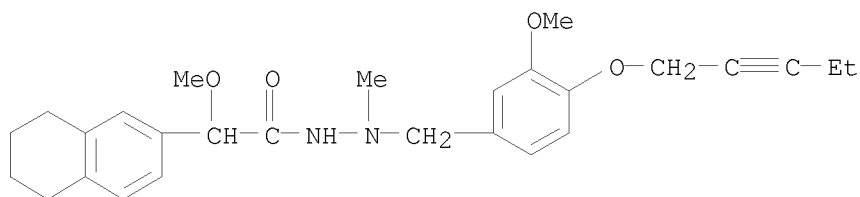
RN 1055788-22-1 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methoxy]- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055788-23-2 CAPLUS

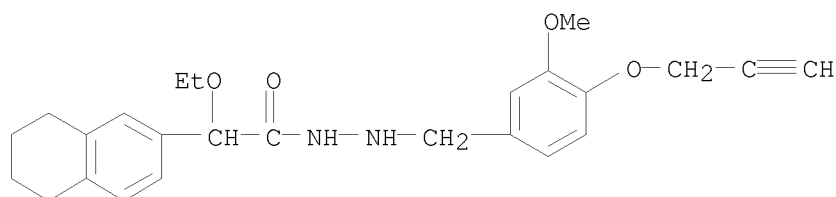
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-,
2-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methyl]-2-methylhydrazide (CA
INDEX NAME)



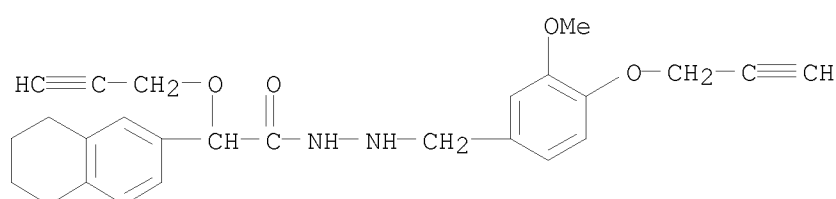
RN 1055788-39-0 CAPLUS

CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
2-[[3-methoxy-4-(2-propyn-1-yloxy)phenyl]methyl]hydrazide (CA INDEX NAME)

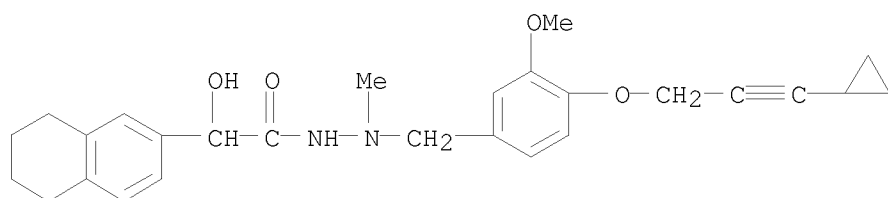
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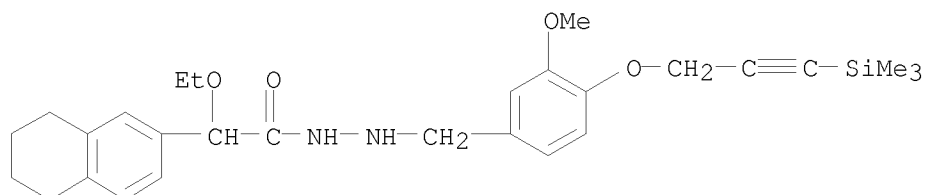
RN 1055788-40-3 CAPLUS
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[[3-methoxy-4-(2-propyn-1-yloxy)phenyl]methyl]hydrazide (CA INDEX NAME)



RN 1055789-27-9 CAPLUS
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-,
2-[[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]methyl]-2-
methylhydrazide (CA INDEX NAME)

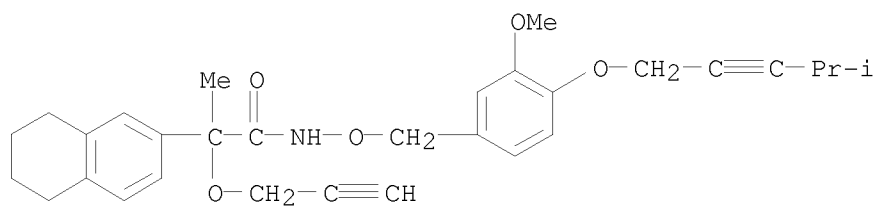


RN 1055789-52-0 CAPLUS
CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
2-[[3-methoxy-4-[[3-(trimethylsilyl)-2-propyn-1-yl]oxy]phenyl]methyl]hydrazide (CA INDEX NAME)

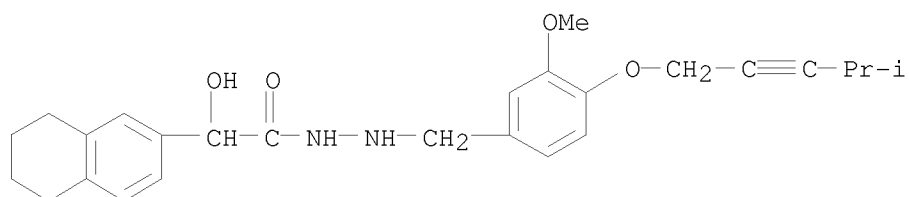


RN 1055789-80-4 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

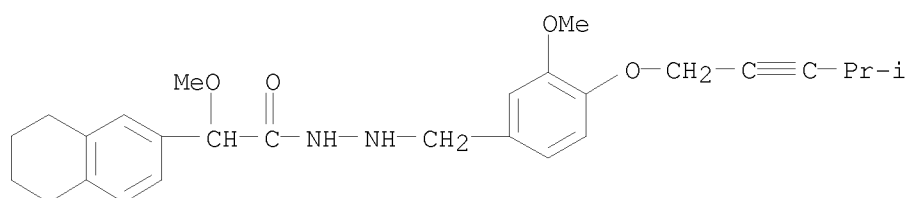
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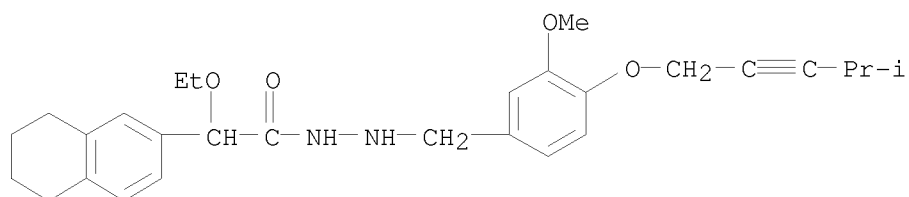
RN 1055789-81-5 CAPLUS
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RN 1055789-82-6 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

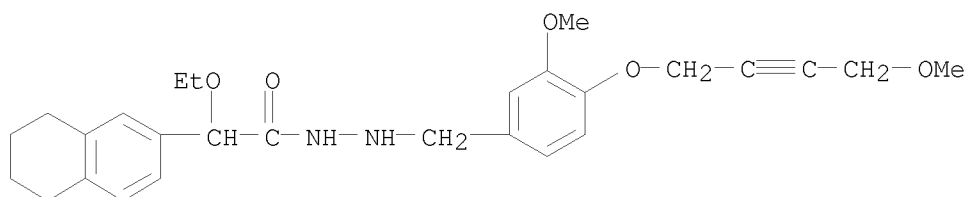


RN 1055789-83-7 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

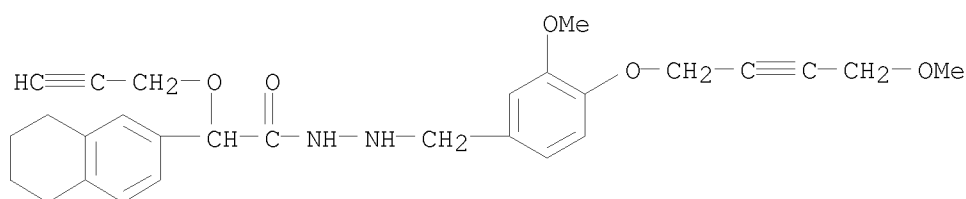


RN 1055790-39-0 CAPLUS
CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
2-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methyl]hydrazide (CA
INDEX NAME)

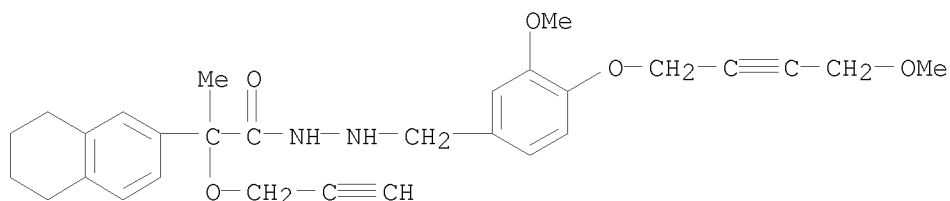
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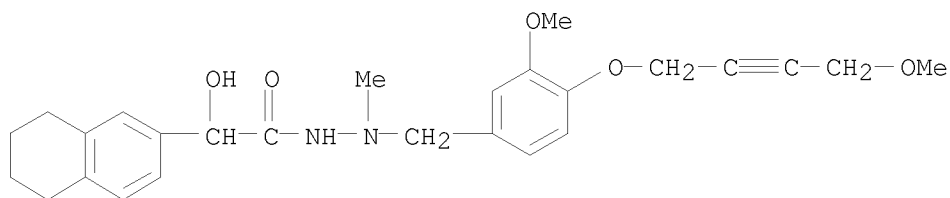
RN 1055790-40-3 CAPLUS
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methyl]hydrazide (CA
INDEX NAME)



RN 1055790-41-4 CAPLUS
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-
propyn-1-yloxy)-, 2-[[3-methoxy-4-[(4-methoxy-2-butyn-1-
yl)oxy]phenyl]methyl]hydrazide (CA INDEX NAME)



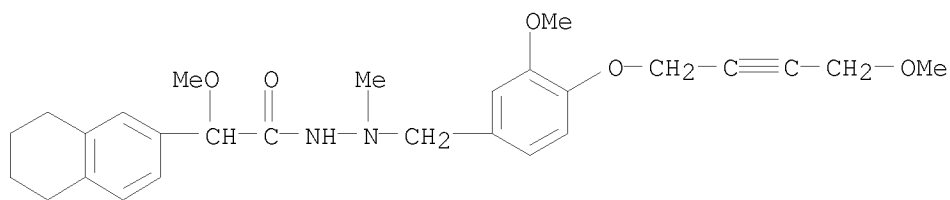
RN 1055790-42-5 CAPLUS
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-,
2-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methyl]-2-
methylhydrazide (CA INDEX NAME)



RN 1055790-43-6 CAPLUS
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-,

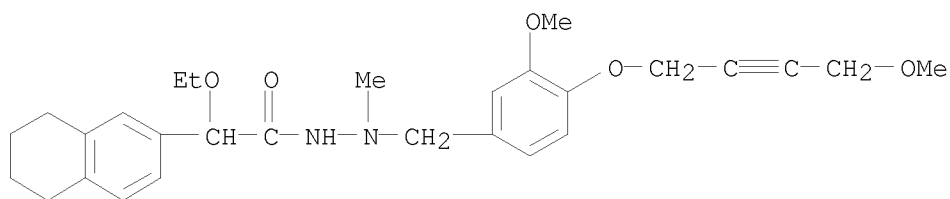
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2-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



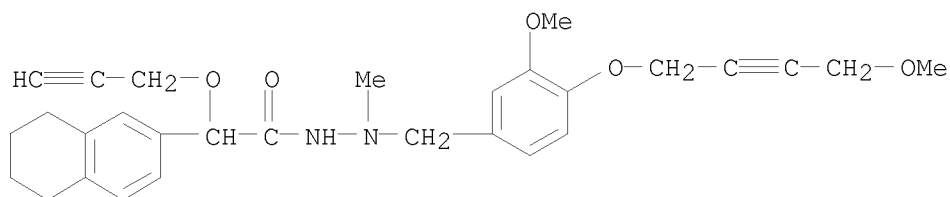
RN 1055790-44-7 CAPLUS

CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-, 2-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



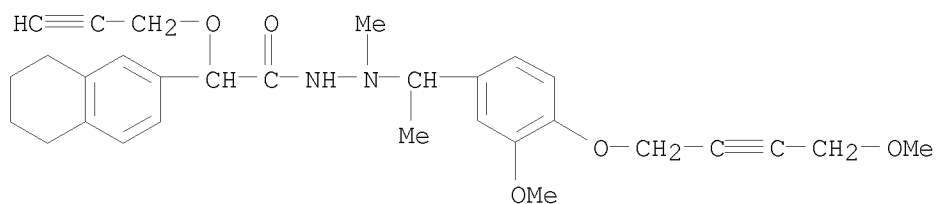
RN 1055790-45-8 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



RN 1055790-46-9 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-, 2-[1-[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]ethyl]-2-methylhydrazide (CA INDEX NAME)



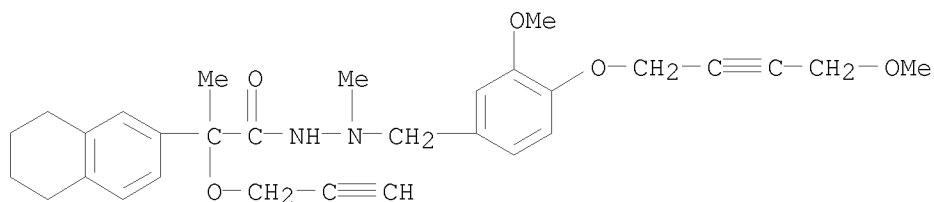
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Erich Leese

10/513699

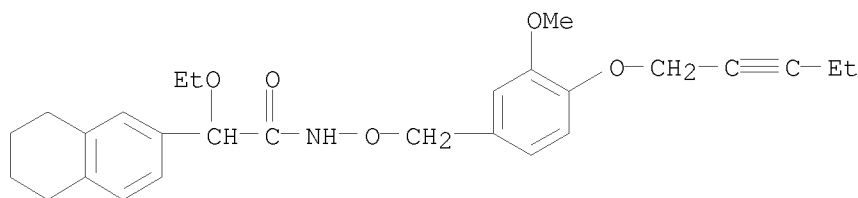
RN 1055790-47-0 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)-, 2-[[3-methoxy-4-[(4-methoxy-2-butyn-1-yl)oxy]phenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



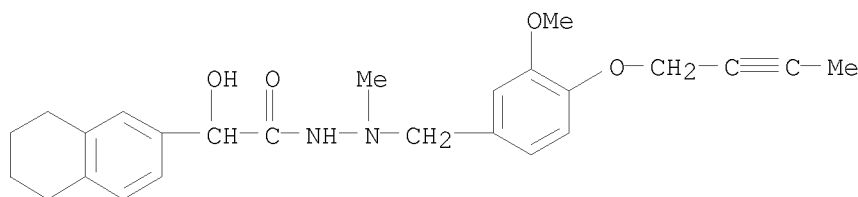
RN 1055792-17-0 CAPLUS

CN 2-Naphthaleneacetamide, α -ethoxy-5,6,7,8-tetrahydro-N-[[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]methoxy]- (CA INDEX NAME)



RN 1055792-28-3 CAPLUS

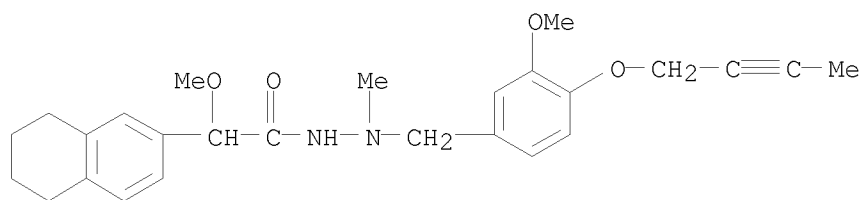
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -hydroxy-, 2-[[4-(2-butyn-1-yloxy)-3-methoxyphenyl]methyl]-2-methylhydrazide (CA INDEX NAME)



RN 1055792-29-4 CAPLUS

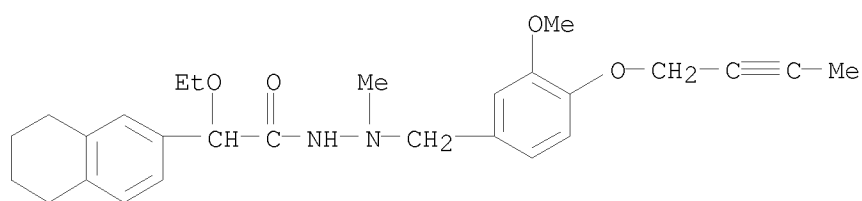
CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-, 2-[[4-(2-butyn-1-yloxy)-3-methoxyphenyl]methyl]-2-methylhydrazide (CA INDEX NAME)

10/513699



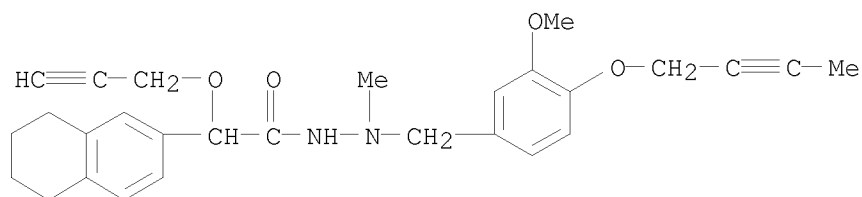
RN 1055792-30-7 CAPLUS

CN 2-Naphthaleneacetic acid, α -ethoxy-5,6,7,8-tetrahydro-,
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INDEX NAME)



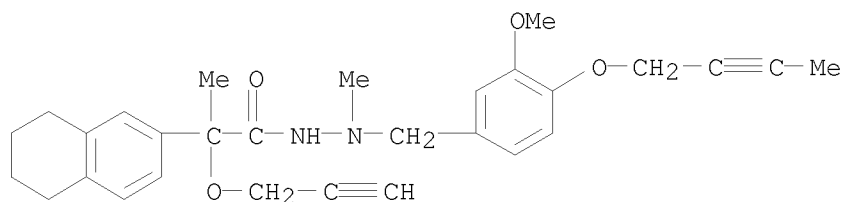
RN 1055792-31-8 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[[4-(2-butyn-1-yloxy)-3-methoxyphenyl]methyl]-2-methylhydrazide (CA
INDEX NAME)



RN 1055792-32-9 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methyl- α -(2-
propyn-1-yloxy)-, 2-[[4-(2-butyn-1-yloxy)-3-methoxyphenyl]methyl]-2-
methylhydrazide (CA INDEX NAME)

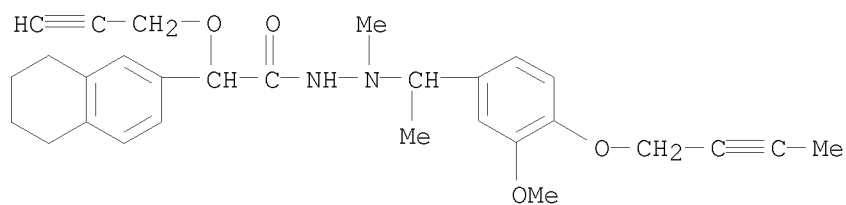


RN 1055792-33-0 CAPLUS

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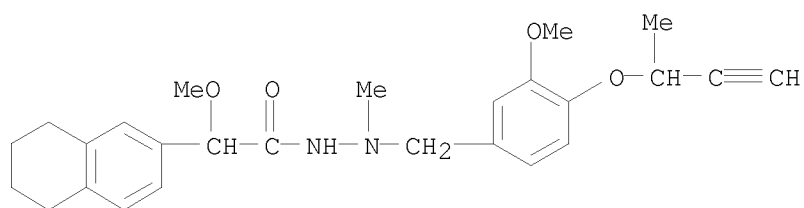
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2-[1-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-2-methylhydrazide (CA
INDEX NAME)



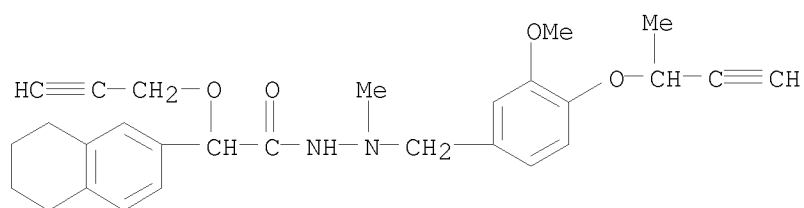
RN 1055792-63-6 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -methoxy-,
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methylhydrazide (CA INDEX NAME)



RN 1055792-66-9 CAPLUS

CN 2-Naphthaleneacetic acid, 5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)-,
2-[[3-methoxy-4-[(1-methyl-2-propyn-1-yl)oxy]phenyl]methyl]-2-
methylhydrazide (CA INDEX NAME)



REFERENCE COUNT:

5

THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L5 ANSWER 4 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2003:1014711 CAPLUS

DOCUMENT NUMBER: 139:403260

TITLE: Preparation of ureidoalkylpiperidines as modulators of chemokine CCR3 receptor activity.

INVENTOR(S): Ko, Soo S.; Delucca, George V.; Duncia, John V.; Santella, Joseph B., III; Wacker, Dean A.

PATENT ASSIGNEE(S): Bristol-Myers Squibb Pharma Company, USA

SOURCE: U.S., 145 pp., Cont.-in-part of U.S. Ser. No. 465,286, abandoned.

CODEN: USXXAM

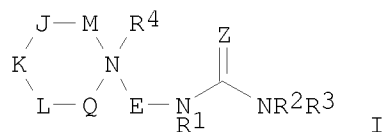
DOCUMENT TYPE: Patent

LANGUAGE: English

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 6605623 B1		20030812US	2000-XI598821 20000621	
PRIORITY APPLN. INFO.:			US 1998-112717P	19981218
			US 1999-161243P	19991022
			US 1999-465286	19991217

GI



AB [Title compds. I; M = CH₂, CHR₅, CHR₁₃, CR₁₃R₁₃, CR₅R₁₃; Q = CH₂, CHR₅, CHR₁₃, CR₁₃R₁₃, CR₅R₁₃; J, L = CH₂, CHR₅, CHR₆, CR₆R₆, CR₅R₆; Z = O, S; M = CH₂, CHR₅, CHR₁₃, CR₁₃R₁₃, CR₅R₁₃; K = CHR₅, CR₅R₆; Z = O, S; E = (CHR₇)(CHR₉)v(CR₁₁R₁₂); R₁, R₂ = H, alkyl, alkenyl, alkynyl, (substituted) alkylcycloalkyl; R₂R₃ = atoms to form a (substituted) 5-7 membered ring; R₃, R₅ = (substituted) (alkyl)cycloalkyl, (alkyl)heterocyclyl; R₄ = null, O, alkyl, alkenyl, alkynyl, etc.; R₄ with R₇, R₉, or R₁₁ = atoms to form a 5-7 membered ring; R₆ = alkyl, alkenyl, alkynyl, etc.; R₇, R₉ = H; R₄R₇, R₄R₉ = (substituted) spirocyclyl; R₁₃ = alkyl, alkenyl, alkynyl, cycloalkyl, etc.; R₁₁R₁₂ = pyrrolidinyl, tetrahydrofuryl, piperidinyl, tetrahydropyranyl; v = 1, 2], were prepared as modulators of chemokine activity (no data) for preventing asthma and other allergic diseases. Thus, 4-benzyl-1-(3-aminopropyl)piperidine (preparation given) in THF was treated with 3-cyanophenyl isocyanate to give N-(3-cyanophenyl)-N'-[3-[4-(phenylmethyl)-1-piperidinyl]propyl]urea. A pharmaceutical composition comprising the compound I was claimed. [This

abstract

record is one of 15 records for this document necessitated by the large number of index entries required to fully index the document and publication system constraints.]

IT 1084141-39-8

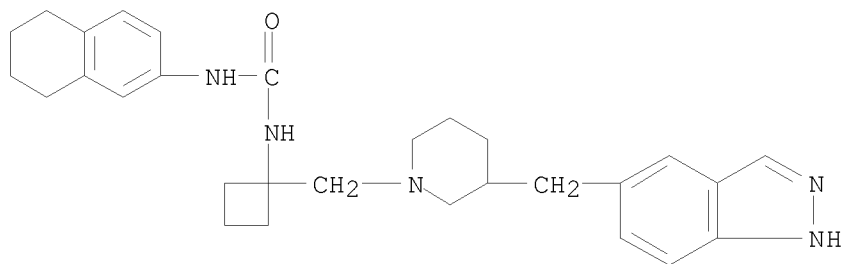
RL: PRPH (Prophetic)

(Preparation of ureidoalkylpiperidines as modulators of chemokine CCR3 receptor activity.)

RN 1084141-39-8 CAPLUS

10/513699

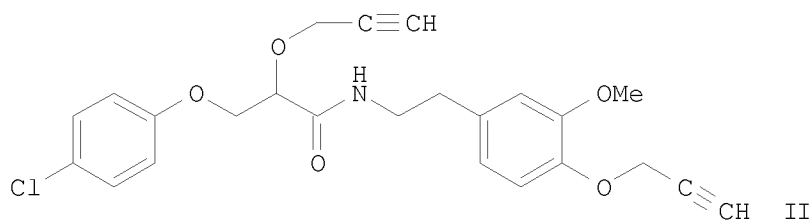
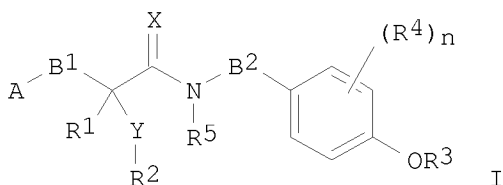
CN Urea, N-[1-[[3-(1H-indazol-5-ylmethyl)-1-piperidinyl]methyl]cyclobutyl]-N'-(5,6,7,8-tetrahydro-2-naphthalenyl)- (CA INDEX NAME)



REFERENCE COUNT: 18 THERE ARE 18 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L5 ANSWER 5 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 2003:1014568 CAPLUS
DOCUMENT NUMBER: 138:411232
TITLE: Preparation of α -oxygenated or α -thiolated
carboxylic acid phenethylamides for controlling fungal
infestation in plants
INVENTOR(S): Zeller, Martin; Lamberth, Clemens; Kriz, Miroslav
PATENT ASSIGNEE(S): Syngenta Participations A.-G., Switz.
SOURCE: PCT Int. Appl., 100 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003042167 A1		20030522	WO 2002-XE12845 20021115	
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZM, ZW				
RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG, TR				
PRIORITY APPLN. INFO.:			GB 2001-27556	20011116
GI				



Erich Leese

heteroarylalkyl, heteroarylalkenyl, or heteroarylalkynyl; R4 = halo, CN, NO2, NH2, CHO, CO2H, (halo)alkyl, (halo)alkenyl(oxy), (halo)alkynyl(oxy), (halo)alkoxy(alkyl), (halo)alkylthio, (halo)alkanoyl, (halo)(di)alkylamino, or (halo)alkoxycarbonyl; R5 = H, alkyl, alkenyl, or alkynyl; n = 0-4; B1 = (CR10R11)q or (CHR10R11)rZ(CR12R13)s; q = 2-4; r = 0-3; s = 1-3; r + s = 1-3; Z = O, S, SO, SO2, NR6, CO, OCO, CO2, NR6CO, or CONR6; R6 = H or alkyl; R10-R13 = independently H or alkyl; B2 = alkylene bridge; and optical isomers and mixts. thereof] were prepared. These compds. possess useful plant protecting properties and may be employed advantageously in agricultural practice for controlling or preventing the infestation of plants by phytopathogenic microorganisms, especially fungi. For example, 3-chlorolactic acid was coupled with 4-chlorophenol in 3.3N NaOH to give 3-(4-chlorophenoxy)-2-hydroxypropionic acid. Amidation with 2-[3-methoxy-4-[(prop-2-ynyl)oxy]phenyl]ethylamine•HCl in the presence of N,N-diisopropylethylamine in DMF, followed by etherification with propargyl bromide in toluene provided the N-(phenethyl)- α -(propargyloxy)propionamide II. The latter showed residual protective action and residual curative action against fungal infestation by *Plasmopara viticola* on vines, *Phytophthora* on tomato plants, and *Phytophthora* on potato plants by 80-100% at 200 ppm. [This abstract record is one of 6 records for this document necessitated by the large number of index entries required to fully index the document and publication system constraints.]

IT	1067826-29-2	1067826-30-5	1067826-31-6
	1067826-32-7	1067826-33-8	1067826-34-9
	1067826-35-0	1067826-36-1	1067826-37-2
	1067835-77-1	1067835-79-3	1067835-80-6
	1067835-81-7	1067835-82-8	1067835-84-0
	1067845-41-3	1067845-42-4	1067845-43-5
	1067845-44-6	1067845-45-7	1067845-46-8
	1067845-47-9	1067845-48-0	1067848-42-3
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	1067873-59-9	1068186-43-5	1068186-44-6
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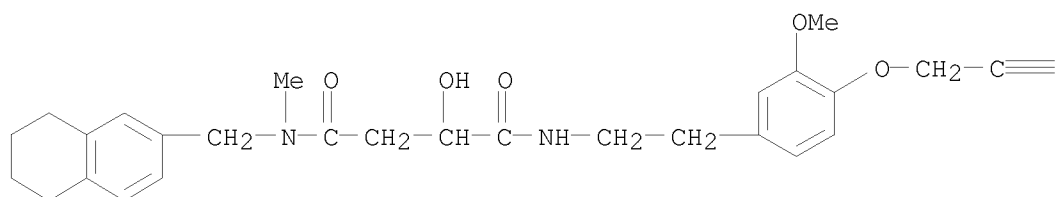
RL: PRPH (Prophetic)

(Preparation of α -oxygenated or α -thiolated carboxylic acid phenethylamides for controlling fungal infestation in plants)

RN 1067826-29-2 CAPLUS

CN Butanediamide, 2-hydroxy-N1-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-N4-methyl-N4-[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]- (CA INDEX NAME)

PAGE 1-A



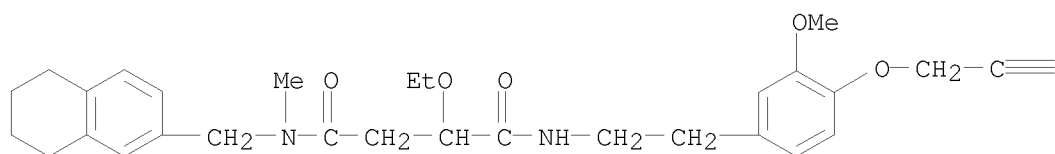
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PAGE 1-A



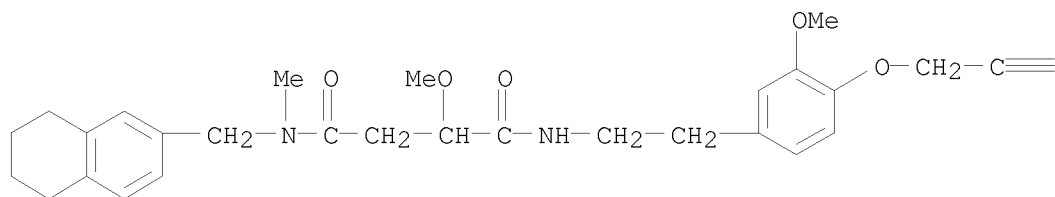
PAGE 1-B

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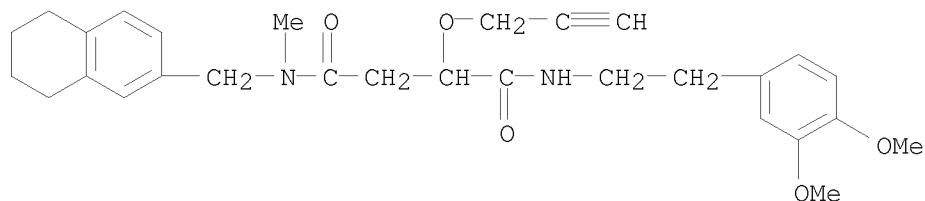
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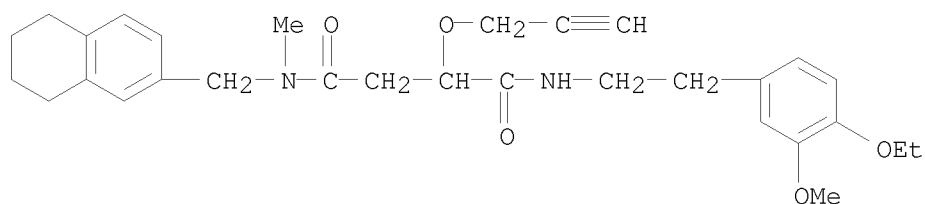
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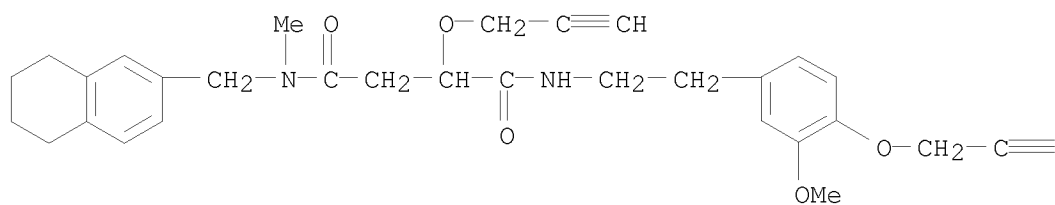
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CN INDEX NAME NOT YET ASSIGNED



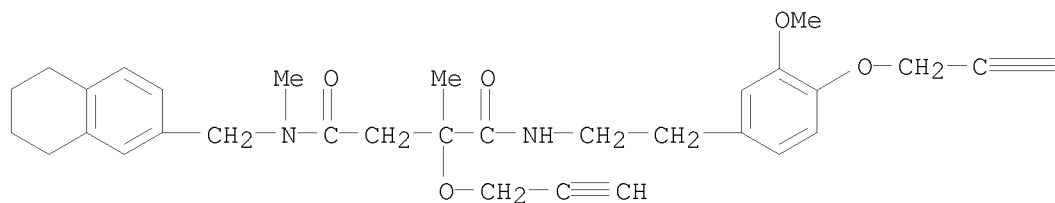
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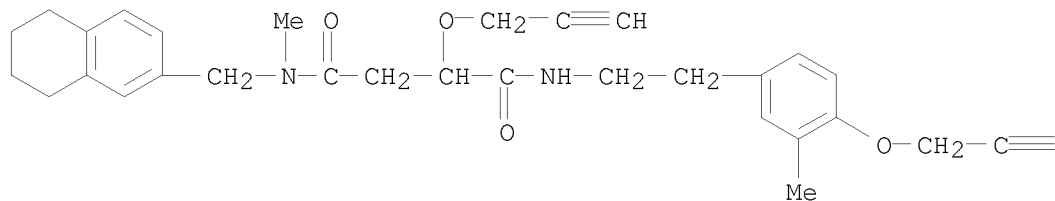


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RN 1067826-36-1 CAPLUS
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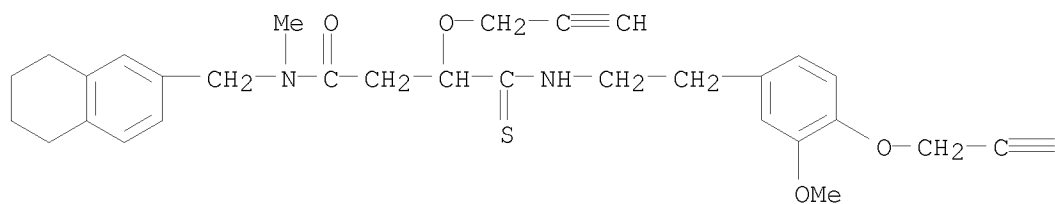


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RN 1067826-37-2 CAPLUS

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PAGE 1-A



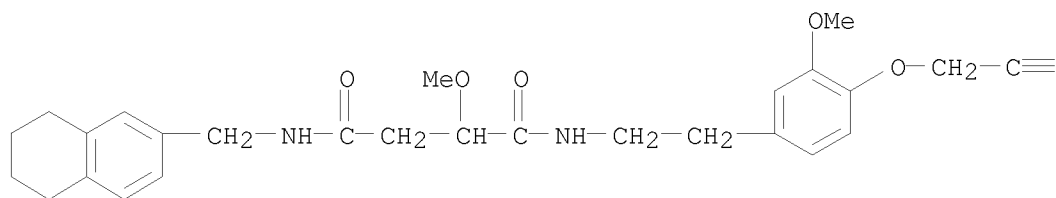
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RN 1067835-77-1 CAPLUS

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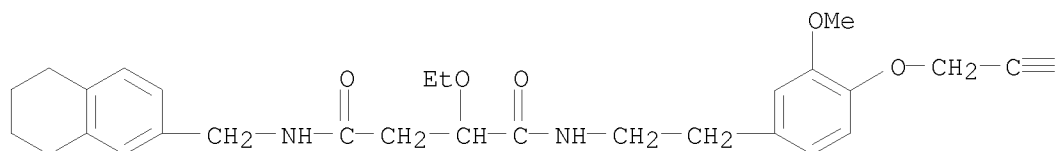
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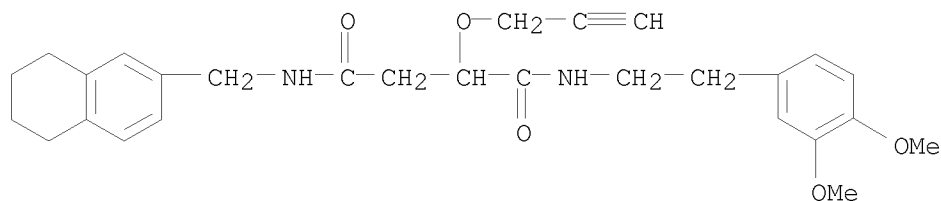
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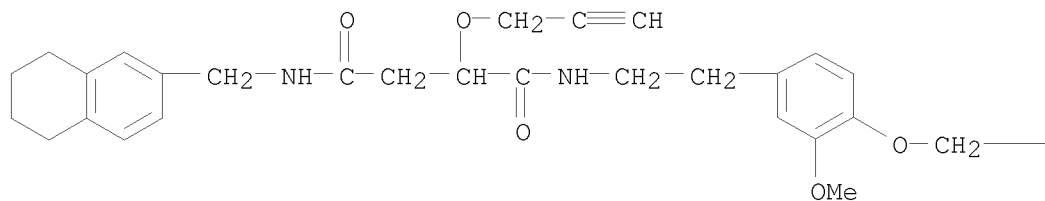
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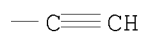
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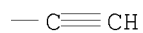
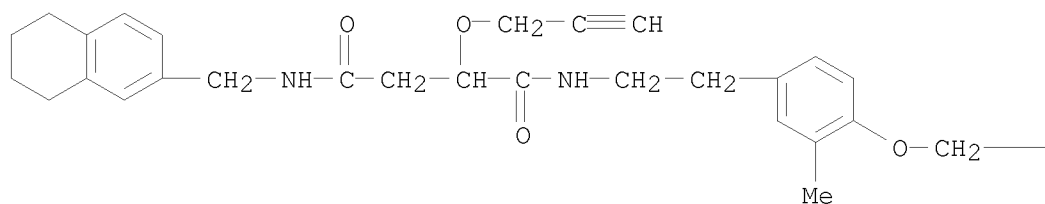
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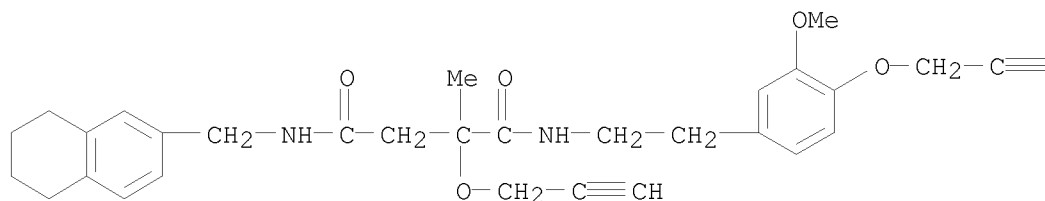
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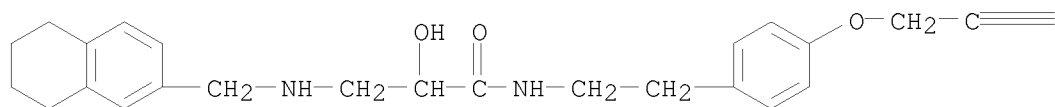
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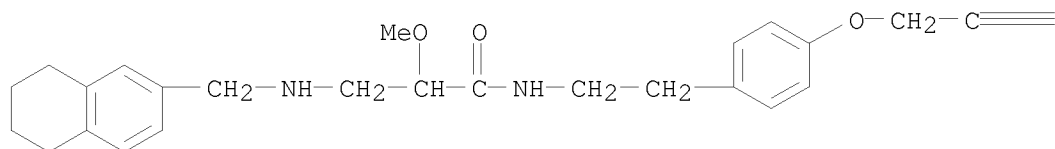


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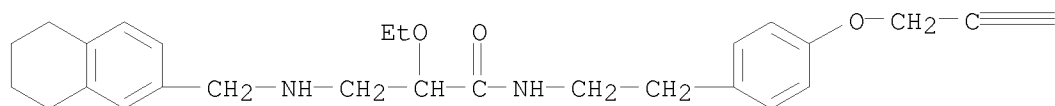
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RN 1067845-42-4 CAPLUS

CN Propanamide, 2-methoxy-N-[2-[4-(2-propyn-1-yloxy)phenyl]ethyl]-3-
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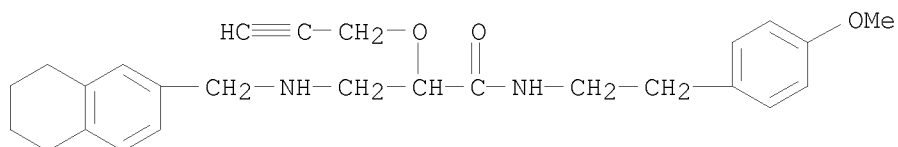
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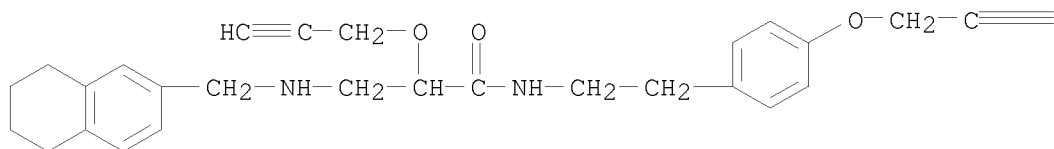


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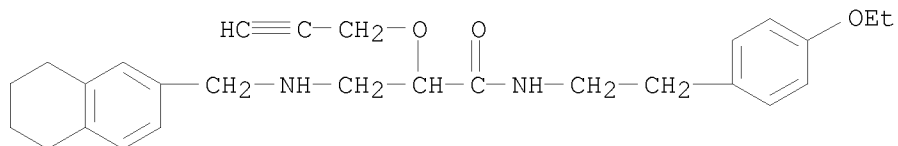
RN 1067845-44-6 CAPLUS
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RN 1067845-45-7 CAPLUS
 CN Propanamide, 2-(2-propyn-1-yloxy)-N-[2-[4-(2-propyn-1-yloxy)phenyl]ethyl]-
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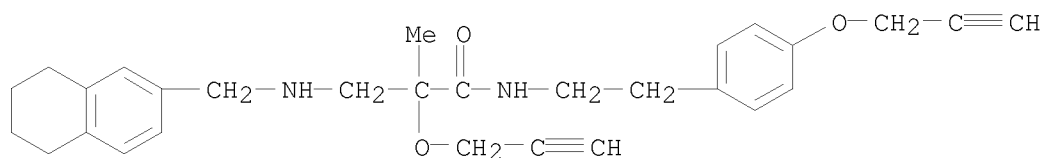
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RN 1067845-46-8 CAPLUS
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RN 1067845-47-9 CAPLUS
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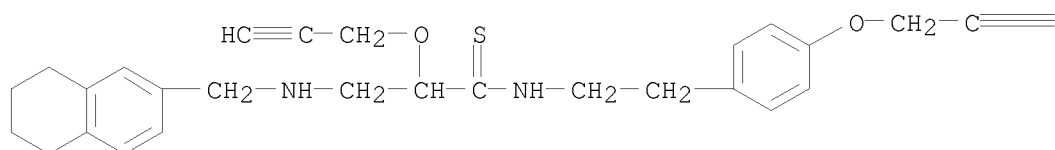
10/513699



RN 1067845-48-0 CAPLUS

CN Propanethioamide, 2-(2-propyn-1-yloxy)-N-[2-[4-(2-propyn-1-yloxy)phenyl]ethyl]-3-[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]amino]- (CA INDEX NAME)

PAGE 1-A



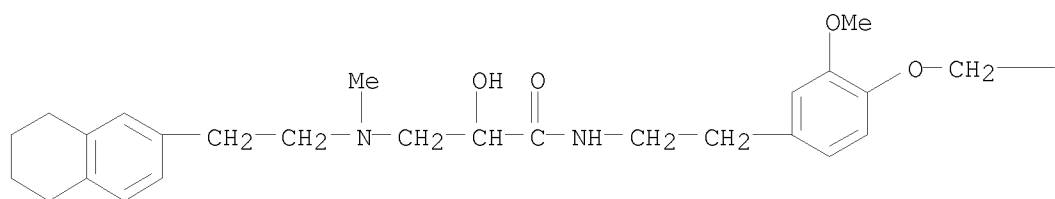
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≡ CH

RN 1067848-42-3 CAPLUS

CN Propanamide, 2-hydroxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-3-[methyl[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)

PAGE 1-A



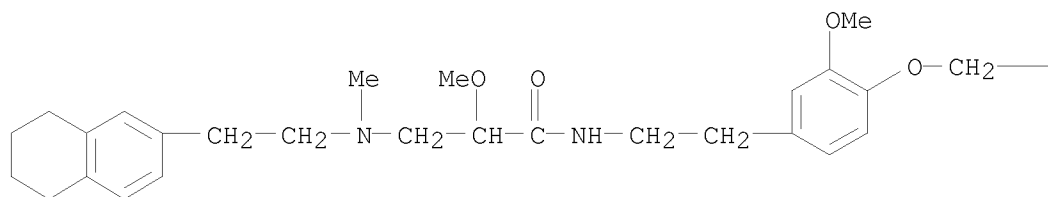
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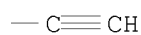
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CN Propanamide, 2-methoxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-3-[methyl[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)

PAGE 1-A

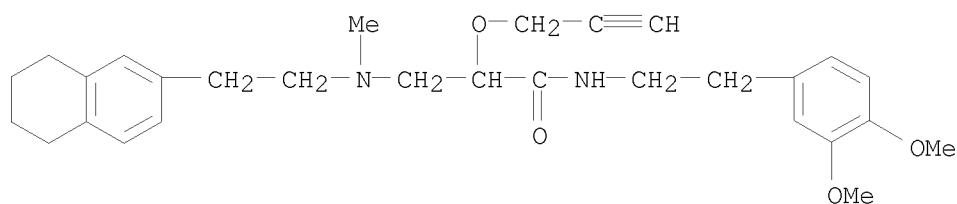


PAGE 1-B



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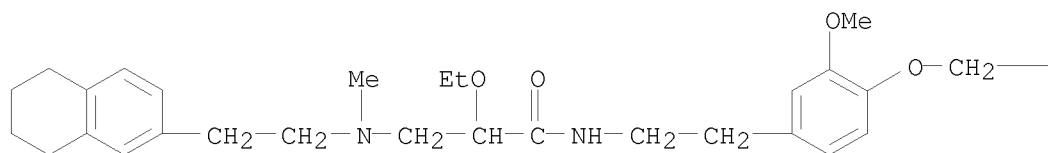
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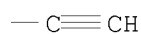
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CN Propanamide, 2-ethoxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-3-[methyl[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)

PAGE 1-A



PAGE 1-B

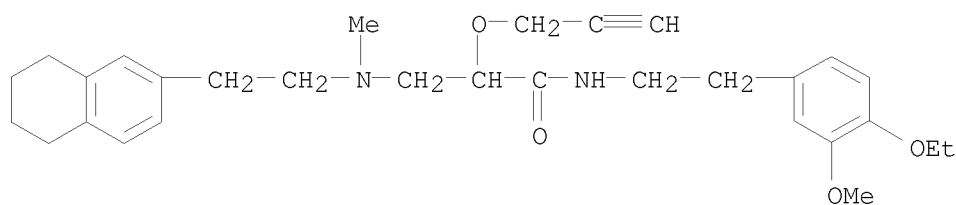


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CN Propanamide, N-[2-(4-ethoxy-3-methoxyphenyl)ethyl]-3-[methyl[2-(5,6,7,8-

10/513699

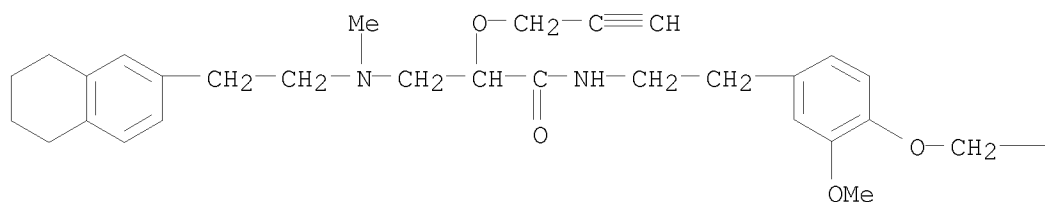
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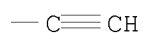
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CN Propanamide, N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-3-[methyl(2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl)amino]-2-(2-propyn-1-yloxy)- (CA INDEX NAME)

PAGE 1-A



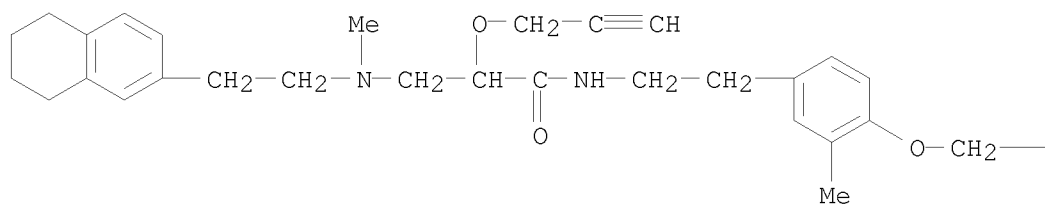
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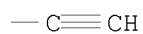
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PAGE 1-A



<12/04/2007>

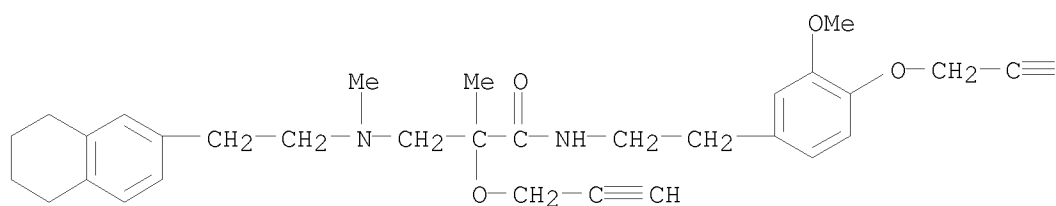
Erich Leese



RN 1067848-49-0 CAPLUS

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PAGE 1-A



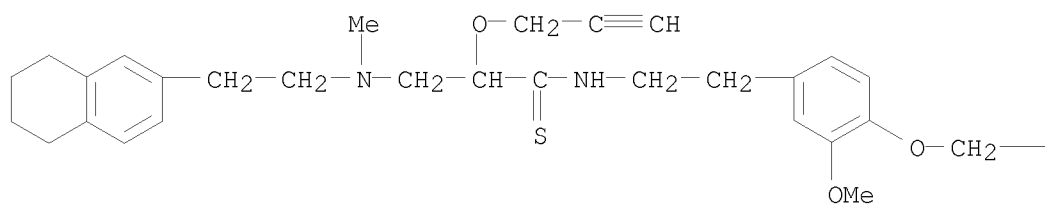
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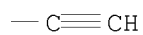
CN Propanethioamide, N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-3-[methyl[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]-2-(2-propyn-1-yloxy)- (CA INDEX NAME)

PAGE 1-A



10/513699

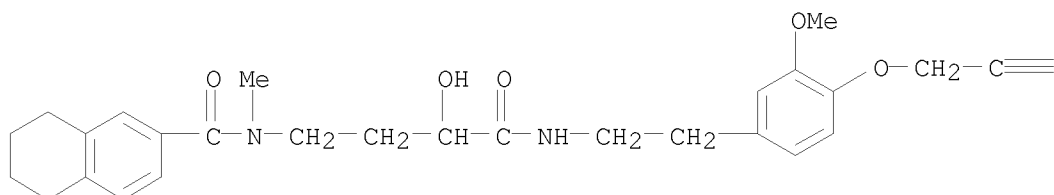
PAGE 1-B



RN 1067859-76-0 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[3-hydroxy-4-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-4-oxobutyl]-N-methyl- (CA INDEX NAME)

PAGE 1-A



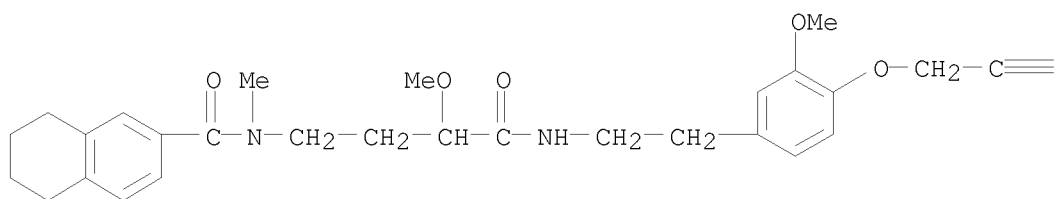
PAGE 1-B



RN 1067859-77-1 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[3-methoxy-4-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-4-oxobutyl]-N-methyl- (CA INDEX NAME)

PAGE 1-A



PAGE 1-B



RN 1067859-78-2 CAPLUS

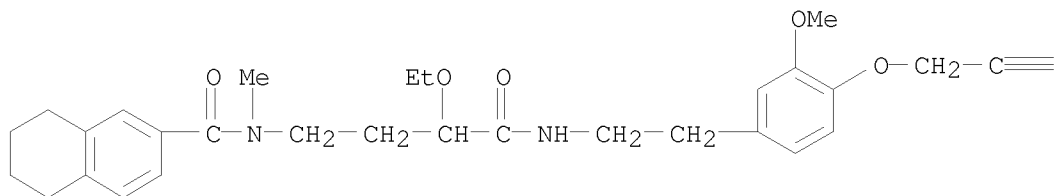
<12/04/2007>

Erich Leese

10/513699

CN 2-Naphthalenecarboxamide, N-[3-ethoxy-4-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-4-oxobutyl]-5,6,7,8-tetrahydro-N-methyl- (CA INDEX NAME)

PAGE 1-A

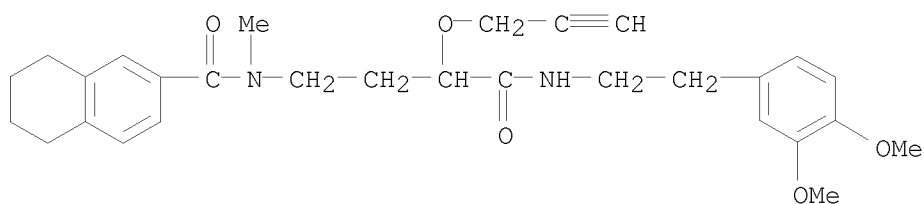


PAGE 1-B

≡CH

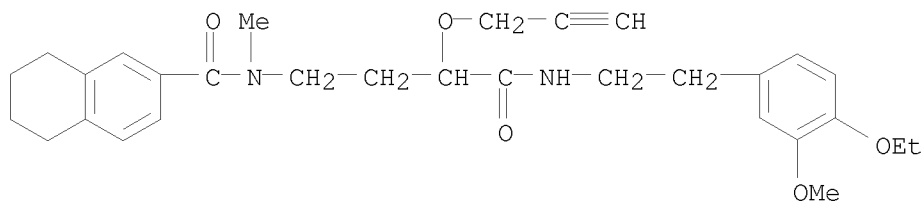
RN 1067859-79-3 CAPLUS

CN 2-Naphthalenecarboxamide, N-[4-[[2-(3,4-dimethoxyphenyl)ethyl]amino]-4-oxo-3-(2-propyn-1-yloxy)butyl]-5,6,7,8-tetrahydro-N-methyl- (CA INDEX NAME)



RN 1067859-80-6 CAPLUS

CN 2-Naphthalenecarboxamide, N-[4-[[2-(4-ethoxy-3-methoxyphenyl)ethyl]amino]-4-oxo-3-(2-propyn-1-yloxy)butyl]-5,6,7,8-tetrahydro-N-methyl- (CA INDEX NAME)



RN 1067859-81-7 CAPLUS

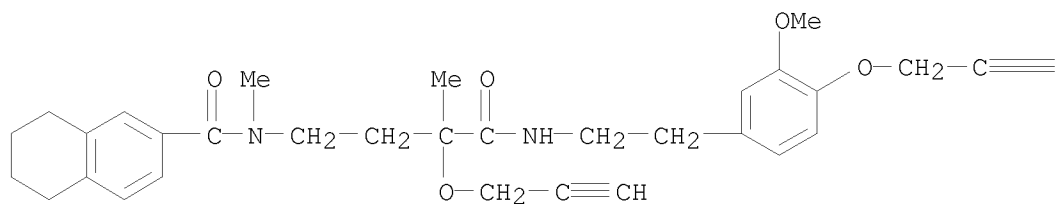
CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[4-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-3-methyl-4-oxo-3-(2-propyn-1-yloxy)butyl]-N-methyl- (CA INDEX NAME)

<12/04/2007>

Erich Leese

10/513699

PAGE 1-A



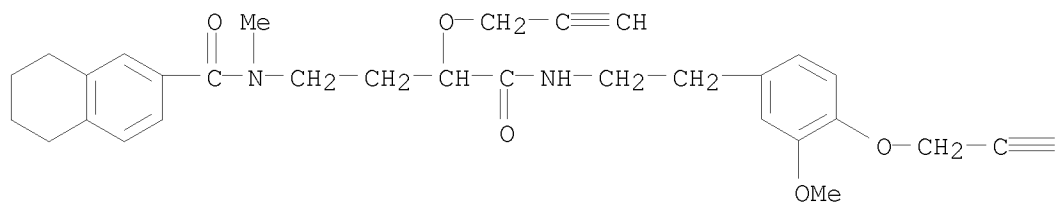
PAGE 1-B

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RN 1067859-82-8 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[4-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-4-oxo-3-(2-propyn-1-yloxy)butyl]-N-methyl- (CA INDEX NAME)

PAGE 1-A



PAGE 1-B

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RN 1067859-83-9 CAPLUS

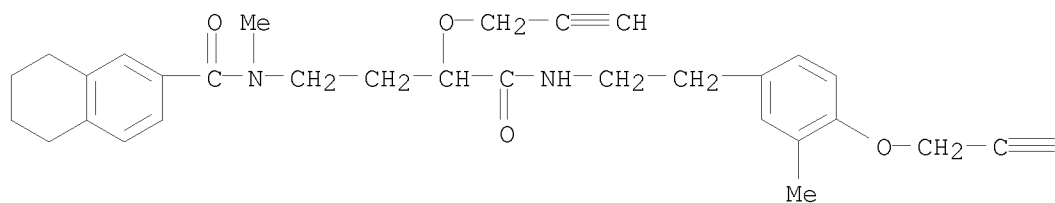
CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-methyl-N-[4-[[2-[3-methyl-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-4-oxo-3-(2-propyn-1-yloxy)butyl]-N-methyl- (CA INDEX NAME)

<12/04/2007>

Erich Leese

10/513699

PAGE 1-A



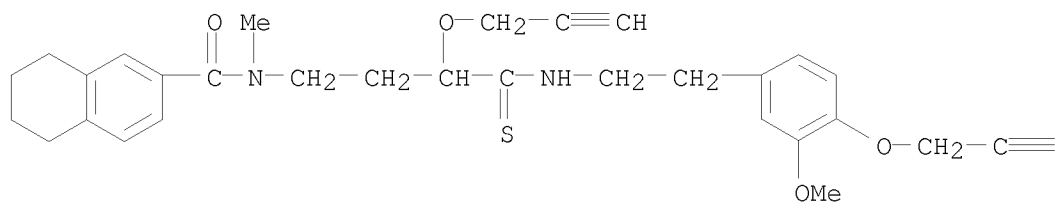
PAGE 1-B

≡ CH

RN 1067859-84-0 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[4-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-3-(2-propyn-1-yloxy)-4-thioxobutyl]-N-methyl- (CA INDEX NAME)

PAGE 1-A



PAGE 1-B

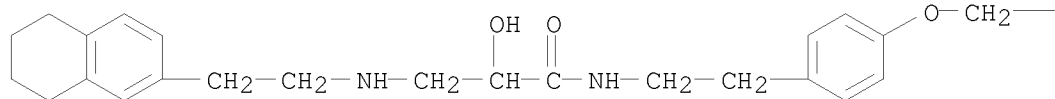
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RN 1067873-52-2 CAPLUS

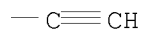
CN Propanamide, 2-hydroxy-N-[2-[4-(2-propyn-1-yloxy)phenyl]ethyl]-3-[[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)

10/513699

PAGE 1-A



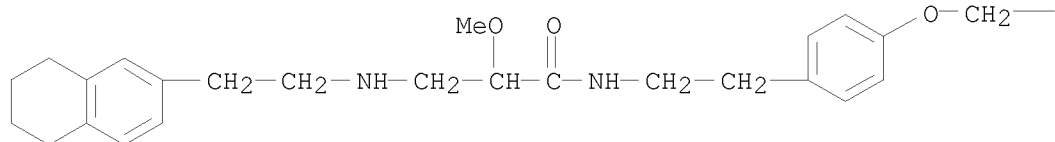
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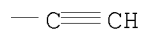
RN 1067873-53-3 CAPLUS

CN Propanamide, 2-methoxy-N-[2-[4-(2-propyn-1-yloxy)phenyl]ethyl]-3-[[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)

PAGE 1-A



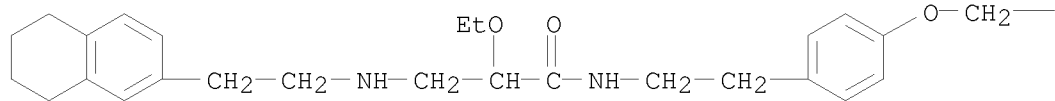
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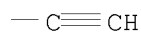
RN 1067873-54-4 CAPLUS

CN Propanamide, 2-ethoxy-N-[2-[4-(2-propyn-1-yloxy)phenyl]ethyl]-3-[[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)

PAGE 1-A



PAGE 1-B



RN 1067873-55-5 CAPLUS

CN Propanamide, N-[2-(4-methoxyphenyl)ethyl]-2-(2-propyn-1-yloxy)-3-[[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)

COc1ccc(cc1)CCNC(=O)C(C#N)CCNCCc2ccc3ccccc3c2CCOC1=CC=CC=C1CCNC(=O)C(C#N)CCNCCc2ccc3ccccc3c2

PAGE 1-A

Chemical structure of a polymer repeat unit:

$$\text{Naphthalene-CH}_2\text{-CH}_2\text{-NH-CH}_2\text{-CH(O-C}\equiv\text{CH)-C(=O)-NH-CH}_2\text{-CH}_2\text{-Ph-O-CH}_2\text{-}$$

The structure shows a naphthalene ring system connected to a polymer chain. The chain consists of a methylene group, a methylene group, an amide group, a methylene group, a chiral center (with a hydroxyl group and a propargyl group), a carbonyl group, an amide group, a methylene group, a methylene group, and a para-substituted phenyl ring. The phenyl ring is connected to a methylene group, which is connected to an oxygen atom. The entire chain is enclosed in brackets with a subscript 'n'.

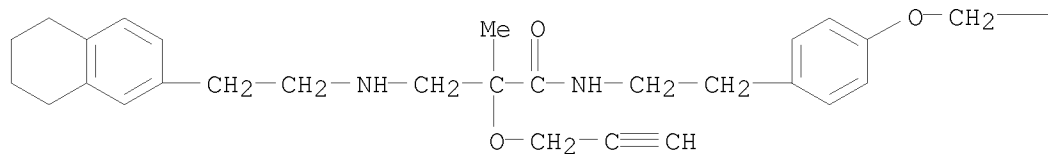
$$-\text{C}\equiv\text{CH}$$

RN	1067873-58-8	CAPLUS
CN	Propanethioamide, 2-(2-propyn-1-yloxy)-N-[2-[4-(2-propyn-1-yloxy)phenyl]ethyl]-3-[[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)	

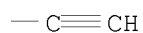
O=C(SCC1=CC=CC=C1OCC2=CC=CC=C2)CNC3=CC=CC=C3OCC4=CC=CC=C4OCC5=CC=CC=C5OCC6=CC=CC=C6OCC7=CC=CC=C7OCC8=CC=CC=C8OCC9=CC=CC=C9OCC10=CC=CC=C10OCC11=CC=CC=C11OCC12=CC=CC=C12OCC13=CC=CC=C13OCC14=CC=CC=C14OCC15=CC=CC=C15OCC16=CC=CC=C16OCC17=CC=CC=C17OCC18=CC=CC=C18OCC19=CC=CC=C19OCC20=CC=CC=C20OCC21=CC=CC=C21OCC22=CC=CC=C22OCC23=CC=CC=C23OCC24=CC=CC=C24OCC25=CC=CC=C25OCC26=CC=CC=C26OCC27=CC=CC=C27OCC28=CC=CC=C28OCC29=CC=CC=C29OCC30=CC=CC=C30OCC31=CC=CC=C31OCC32=CC=CC=C32OCC33=CC=CC=C33OCC34=CC=CC=C34OCC35=CC=CC=C35OCC36=CC=CC=C36OCC37=CC=CC=C37OCC38=CC=CC=C38OCC39=CC=CC=C39OCC40=CC=CC=C40OCC41=CC=CC=C41OCC42=CC=CC=C42OCC43=CC=CC=C43OCC44=CC=CC=C44OCC45=CC=CC=C45OCC46=CC=CC=C46OCC47=CC=CC=C47OCC48=CC=CC=C48OCC49=CC=CC=C49OCC50=CC=CC=C50OCC51=CC=CC=C51OCC52=CC=CC=C52OCC53=CC=CC=C53OCC54=CC=CC=C54OCC55=CC=CC=C55OCC56=CC=CC=C56OCC57=CC=CC=C57OCC58=CC=CC=C58OCC59=CC=CC=C59OCC60=CC=CC=C60OCC61=CC=CC=C61OCC62=CC=CC=C62OCC63=CC=CC=C63OCC64=CC=CC=C64OCC65=CC=CC=C65OCC66=CC=CC=C66OCC67=CC=CC=C67OCC68=CC=CC=C68OCC69=CC=CC=C69OCC70=CC=CC=C70OCC71=CC=CC=C71OCC72=CC=CC=C72OCC73=CC=CC=C73OCC74=CC=CC=C74OCC75=CC=CC=C75OCC76=CC=CC=C76OCC77=CC=CC=C77OCC78=CC=CC=C78OCC79=CC=CC=C79OCC80=CC=CC=C80OCC81=CC=CC=C81OCC82=CC=CC=C82OCC83=CC=CC=C83OCC84=CC=CC=C84OCC85=CC=CC=C85OCC86=CC=CC=C86OCC87=CC=CC=C87OCC88=CC=CC=C88OCC89=CC=CC=C89OCC90=CC=CC=C90OCC91=CC=CC=C91OCC92=CC=CC=C92OCC93=CC=CC=C93OCC94=CC=CC=C94OCC95=CC=CC=C95OCC96=CC=CC=C96OCC97=CC=CC=C97OCC98=CC=CC=C98OCC99=CC=CC=C99OCC100=CC=CC=C100OCC101=CC=CC=C101OCC102=CC=CC=C102OCC103=CC=CC=C103OCC104=CC=CC=C104OCC105=CC=CC=C105OCC106=CC=CC=C106OCC107=CC=CC=C107OCC108=CC=CC=C108OCC109=CC=CC=C109OCC110=CC=CC=C110OCC111=CC=CC=C111OCC112=CC=CC=C112OCC113=CC=CC=C113OCC114=CC=CC=C114OCC115=CC=CC=C115OCC116=CC=CC=C116OCC117=CC=CC=C117OCC118=CC=CC=C118OCC119=CC=CC=C119OCC120=CC=CC=C120OCC121=CC=CC=C121OCC122=CC=CC=C122OCC123=CC=CC=C123OCC124=CC=CC=C124OCC125=CC=CC=C125OCC126=CC=CC=C126OCC127=CC=CC=C127OCC128=CC=CC=C128OCC129=CC=CC=C129OCC130=CC=CC=C130OCC131=CC=CC=C131OCC132=CC=CC=C132OCC133=CC=CC=C133OCC134=CC=CC=C134OCC135=CC=CC=C135OCC136=CC=CC=C136OCC137=CC=CC=C137OCC138=CC=CC=C138OCC139=CC=CC=C139OCC140=CC=CC=C140OCC141=CC=CC=C141OCC142=CC=CC=C142OCC143=CC=CC=C143OCC144=CC=CC=C144OCC145=CC=CC=C145OCC146=CC=CC=C146OCC147=CC=CC=C147OCC148=CC=CC=C148OCC149=CC=CC=C149OCC150=CC=CC=C150OCC151=CC=CC=C151OCC152=CC=CC=C152OCC153=CC=CC=C153OCC154=CC=CC=C154OCC155=CC=CC=C155OCC156=CC=CC=C156OCC157=CC=CC=C157OCC158=CC=CC=C158OCC159=CC=CC=C159OCC160=CC=CC=C160OCC161=CC=CC=C161OCC162=CC=CC=C162OCC163=CC=CC=C163OCC164=CC=CC=C164OCC165=CC=CC=C165OCC166=CC=CC=C166OCC167=CC=CC=C167OCC168=CC=CC=C168OCC169=CC=CC=C169OCC170=CC=CC=C170OCC171=CC=CC=C171OCC172=CC=CC=C172OCC173=CC=CC=C173OCC174=CC=CC=C174OCC175=CC=CC=C175OCC176=CC=CC=C176OCC177=CC=CC=C177OCC178=CC=CC=C178OCC179=CC=CC=C179OCC180=CC=CC=C180OCC181=CC=CC=C181OCC182=CC=CC=C182OCC183=CC=CC=C183OCC184=CC=CC=C184OCC185=CC=CC=C185OCC186=CC=CC=C186OCC187=CC=CC=C187OCC188=CC=CC=C188OCC189=CC=CC=C189OCC190=CC=CC=C190OCC191=CC=CC=C191OCC192=CC=CC=C192OCC193=CC=CC=C193OCC194=CC=CC=C194OCC195=CC=CC=C195OCC196=CC=CC=C196OCC197=CC=CC=C197OCC198=CC=CC=C198OCC199=CC=CC=C199OCC200=CC=CC=C200OCC201=CC=CC=C201OCC202=CC=CC=C202OCC203=CC=CC=C203OCC204=CC=CC=C204OCC205=CC=CC=C205OCC206=CC=CC=C206OCC207=CC=CC=C207OCC208=CC=CC=C208OCC209=CC=CC=C209OCC210=CC=CC=C210OCC211=CC=CC=C211OCC212=CC=CC=C212OCC213=CC=CC=C213OCC214=CC=CC=C214OCC215=CC=CC=C215OCC216=CC=CC=C216OCC217=CC=CC=C217OCC218=CC=CC=C218OCC219=CC=CC=C219OCC220=CC=CC=C220OCC221=CC=CC=C221OCC222=CC=CC=C222OCC223=CC=CC=C223OCC224=CC=CC=C224OCC225=CC=CC=C225OCC226=CC=CC=C226OCC227=CC=CC=C227OCC228=CC=CC=C228OCC229=CC=CC=C229OCC230=CC=CC=C230OCC231=CC=CC=C231OCC232=CC=CC=C232OCC233=CC=CC=C233OCC234=CC=CC=C234OCC235=CC=CC=C235OCC236=CC=CC=C236OCC237=CC=CC=C237OCC238=CC=CC=C238OCC239=CC=CC=C239OCC240=CC=CC=C240OCC241=CC=CC=C241OCC242=CC=CC=C242OCC243=CC=CC=C243OCC244=CC=CC=C244OCC245=CC=CC=C245OCC246=CC=CC=C246OCC247=CC=CC=C247OCC248=CC=CC=C248OCC249=CC=CC=C249OCC250=CC=CC=C250OCC251=CC=CC=C251OCC252=CC=CC=C252OCC253=CC=CC=C253OCC254=CC=CC=C254OCC255=CC=CC=C255OCC256=CC=CC=C256OCC257=CC=CC=C257OCC258=CC=CC=C258OCC259=CC=CC=C259OCC260=CC=CC=C260OCC261=CC=CC=C261OCC262=CC=CC=C262OCC263=CC=CC=C263OCC264=CC=CC=C264OCC265=CC=CC=C265OCC266=CC=CC=C266OCC267=CC=CC=C267OCC268=CC=CC=C268OCC269=CC=CC=C269OCC270=CC=CC=C270OCC271=CC=CC=C271OCC272=CC=CC=C272OCC273=CC=CC=C273OCC274=CC=CC=C274OCC275=CC=CC=C275OCC276=CC=CC=C276OCC277=CC=CC=C277OCC278=CC=CC=C278OCC279=CC=CC=C279OCC280=CC=CC=C280OCC281=CC=CC=C281OCC282=CC=CC=C282OCC283=CC=CC=C283OCC284=CC=CC=C284OCC285=CC=CC=C285OCC286=CC=CC=C286OCC287=CC=CC=C287OCC288=CC=CC=C288OCC289=CC=CC=C289OCC290=CC=CC=C290OCC291=CC=CC=C291OCC292=CC=CC=C292OCC293=CC=CC=C293OCC294=CC=CC=C294OCC295=CC=CC=C295OCC296=CC=CC=C296OCC297=CC=CC=C297OCC298=CC=CC=C298OCC299=CC=CC=C299OCC300=CC=CC=C300OCC301=CC=CC=C301OCC302=CC=CC=C302OCC303=CC=CC=C303OCC304=CC=CC=C304OCC305=CC=CC=C305OCC306=CC=CC=C306OCC307=CC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Erich Leese

PAGE 1-A



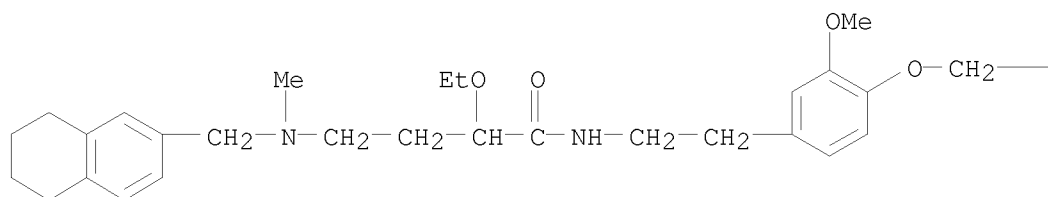
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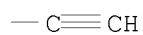
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CN Butanamide, 2-ethoxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-4-[methyl[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]amino]- (CA INDEX NAME)

PAGE 1-A



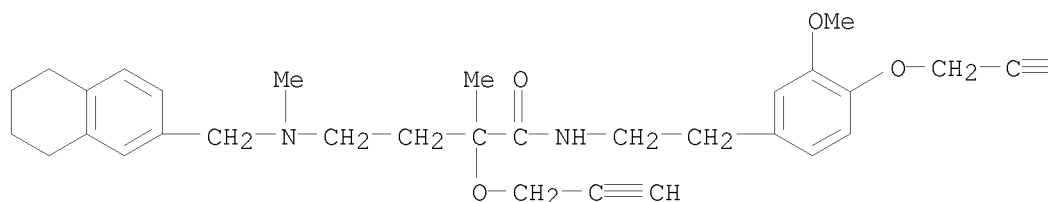
PAGE 1-B



RN 1068186-44-6 CAPLUS

CN Butanamide, N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-2-methyl-4-[methyl[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]amino]-2-(2-propyn-1-yloxy)- (CA INDEX NAME)

PAGE 1-A



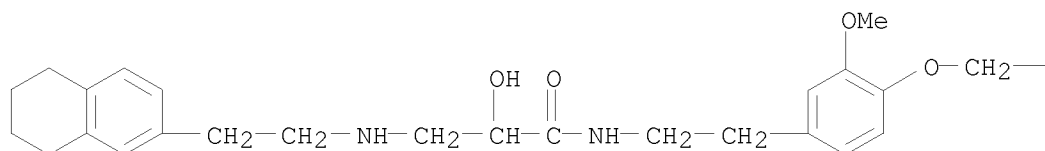
PAGE 1-B



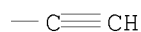
RN 1068190-31-7 CAPLUS

 CN Propanamide, 2-hydroxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-3-
 [[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)

PAGE 1-A



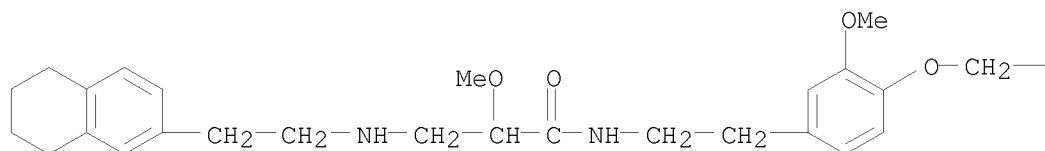
PAGE 1-B



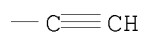
RN 1068190-32-8 CAPLUS

 CN Propanamide, 2-methoxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-3-
 [[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)

PAGE 1-A



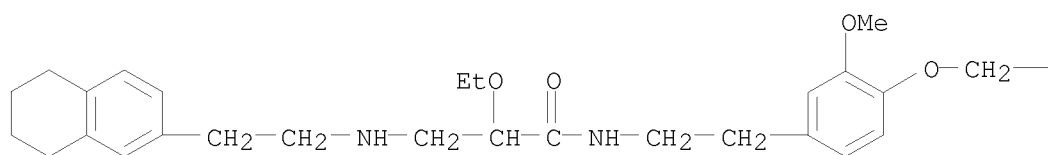
PAGE 1-B



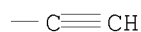
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 CN Propanamide, 2-ethoxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-3-
 [[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)

PAGE 1-A

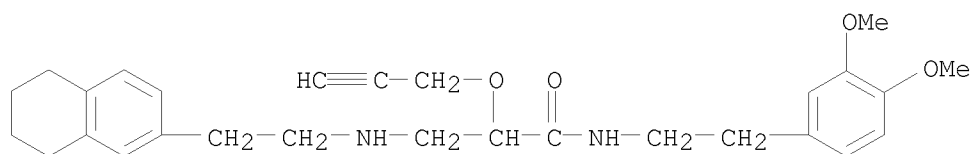


PAGE 1-B



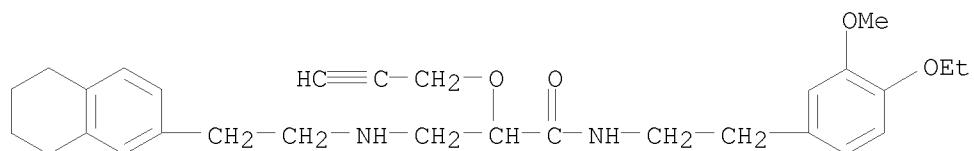
RN 1068190-34-0 CAPLUS

CN Propanamide, N-[2-(3,4-dimethoxyphenyl)ethyl]-2-(2-propyn-1-yloxy)-3-[[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)



RN 1068190-35-1 CAPLUS

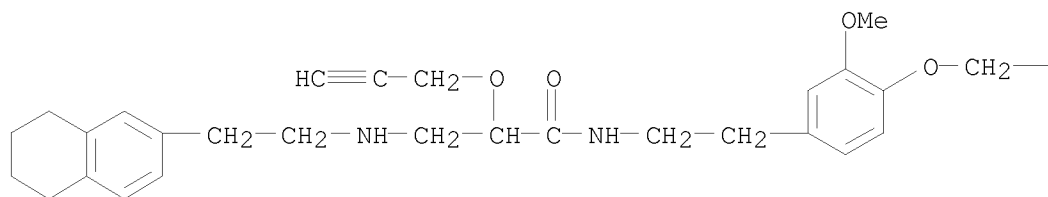
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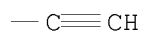
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CN Propanamide, N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-2-(2-propyn-1-yloxy)-3-[[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]- (CA INDEX NAME)

PAGE 1-A



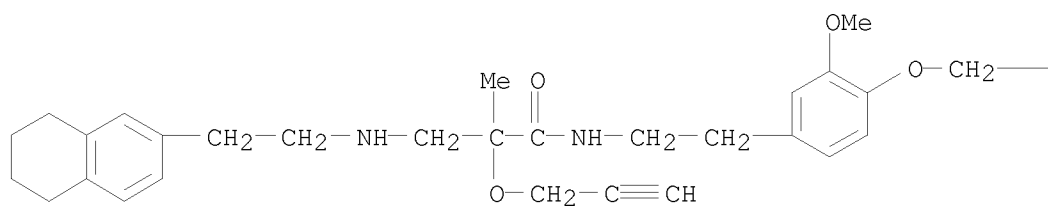
PAGE 1-B



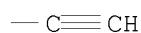
RN 1068190-37-3 CAPLUS

CN Propanamide, N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-2-methyl-2-(2-propyn-1-yloxy)-3-[[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]amino]-(CA INDEX NAME)

PAGE 1-A



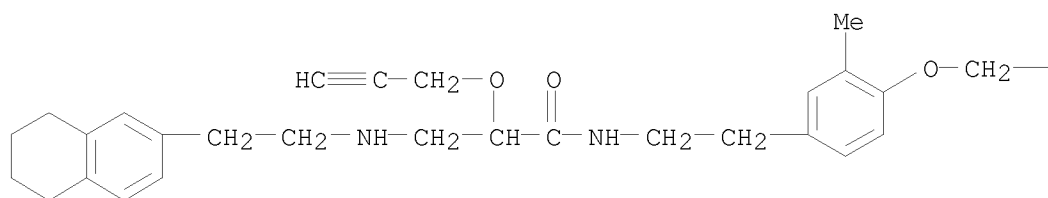
PAGE 1-B



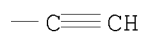
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PAGE 1-A



PAGE 1-B

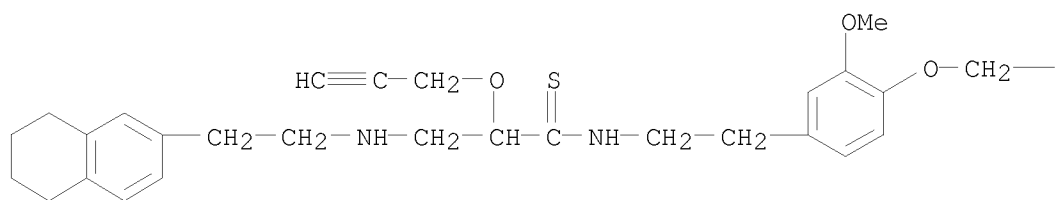


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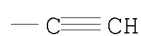
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10/513699

PAGE 1-A

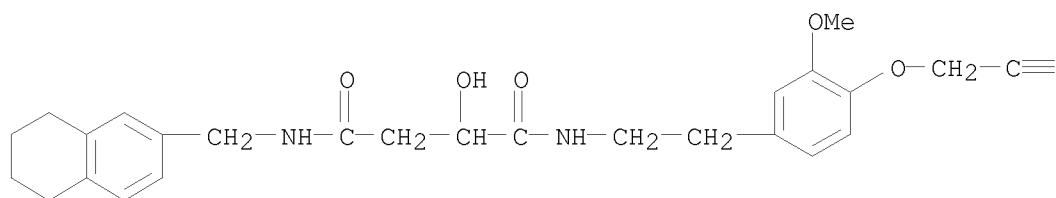


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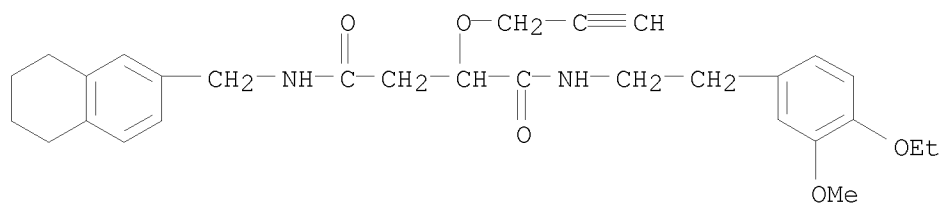
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PAGE 1-B



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CN INDEX NAME NOT YET ASSIGNED



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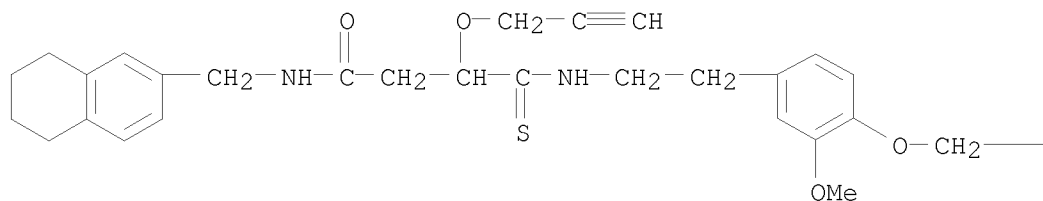
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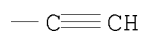
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PAGE 1-A



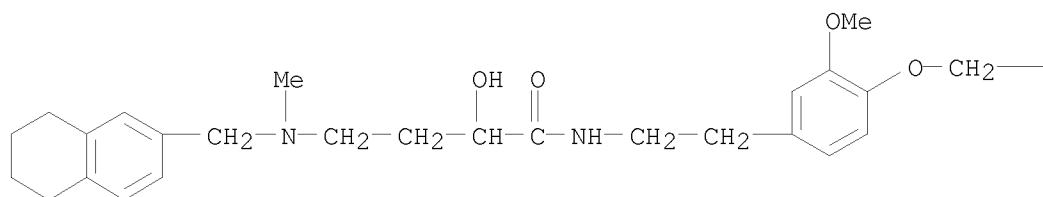
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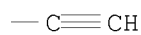
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PAGE 1-A

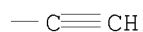
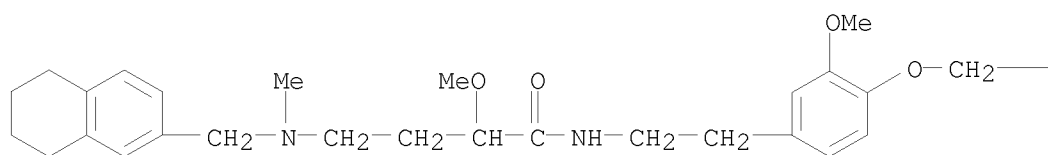


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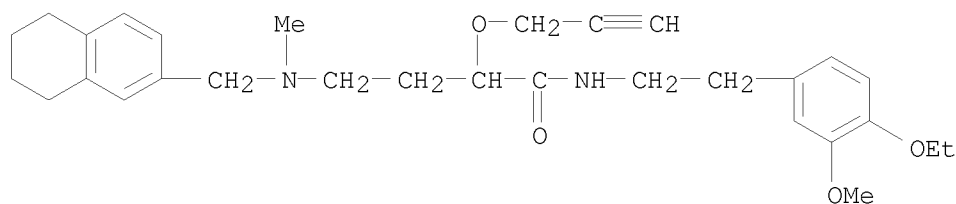


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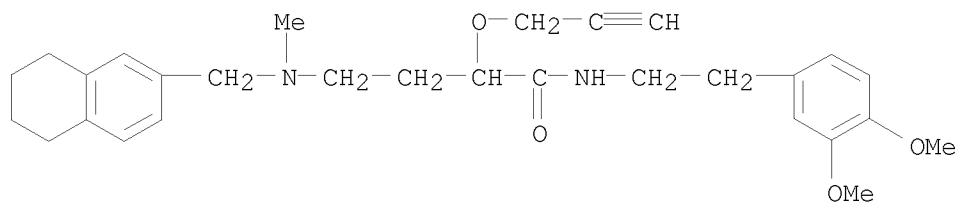
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RN 1068197-83-0 CAPLUS
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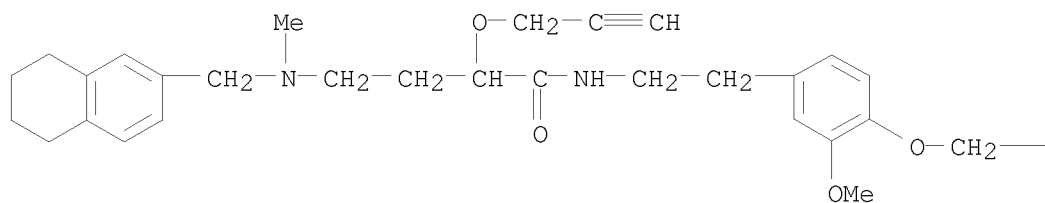
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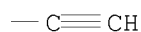
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10/513699

PAGE 1-A

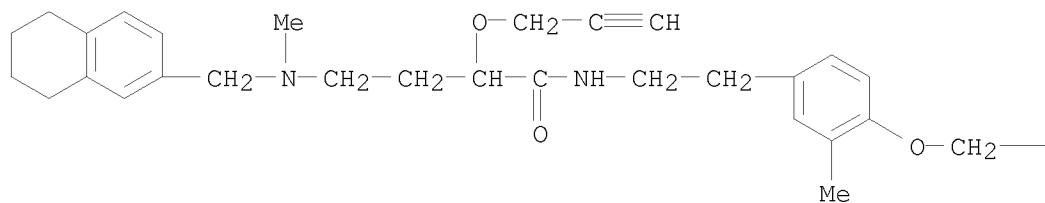


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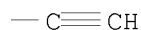


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PAGE 1-A



PAGE 1-B



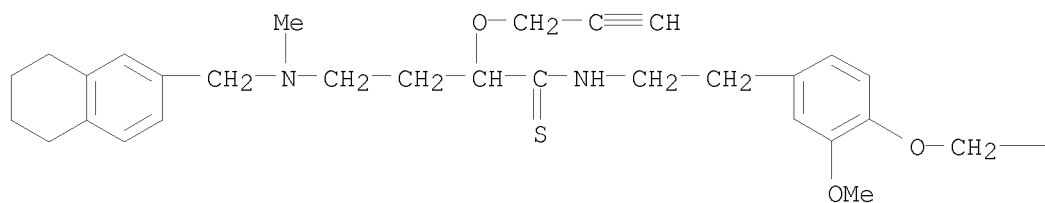
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10/513699

PAGE 1-A

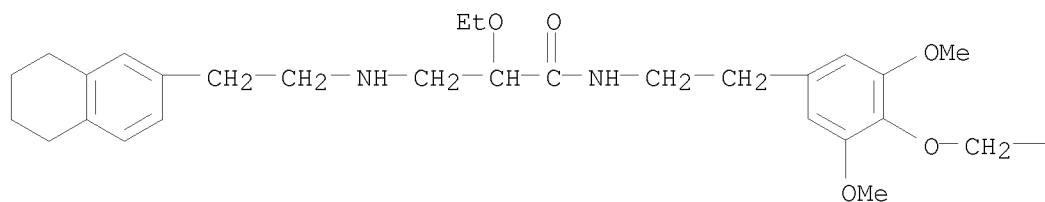


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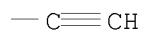


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PAGE 1-A



PAGE 1-B

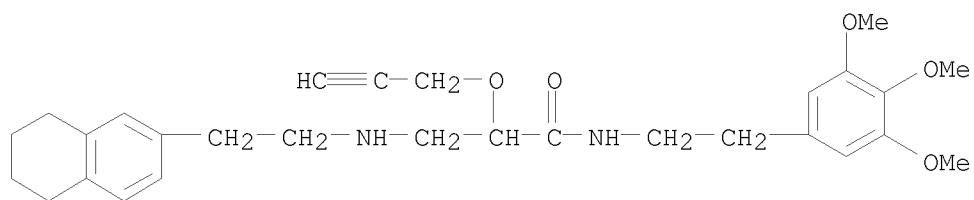


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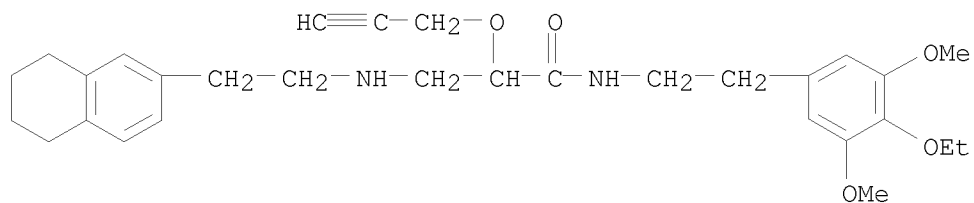
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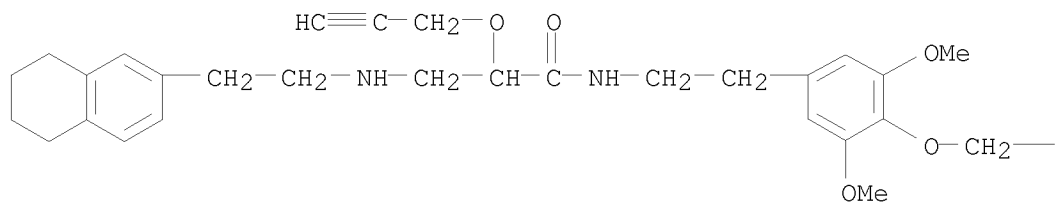


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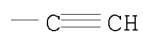


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PAGE 1-A



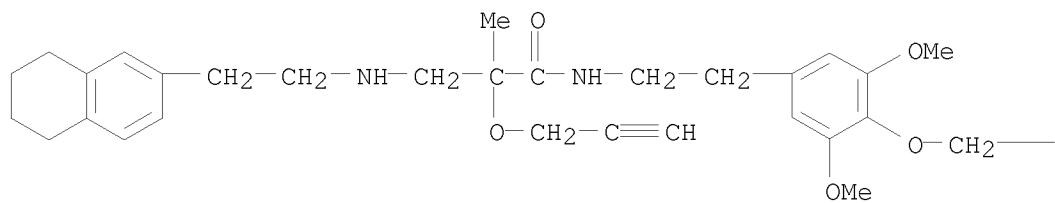
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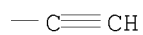
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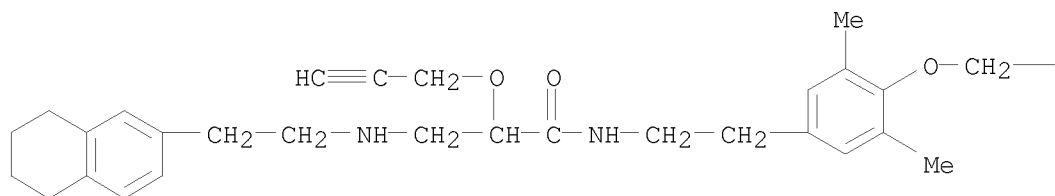


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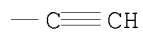


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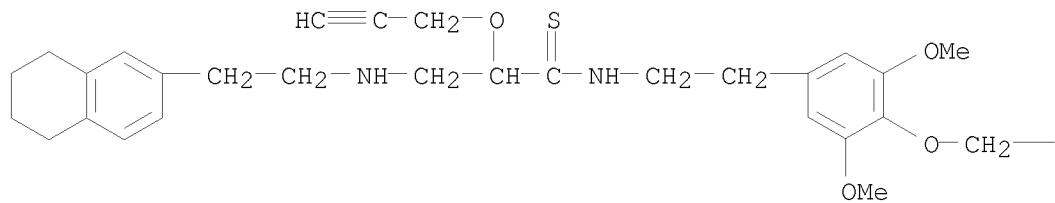


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PAGE 1-A



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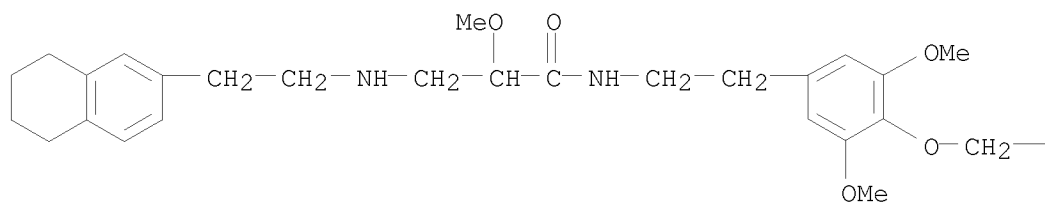
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PAGE 1-B

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PAGE 1-A

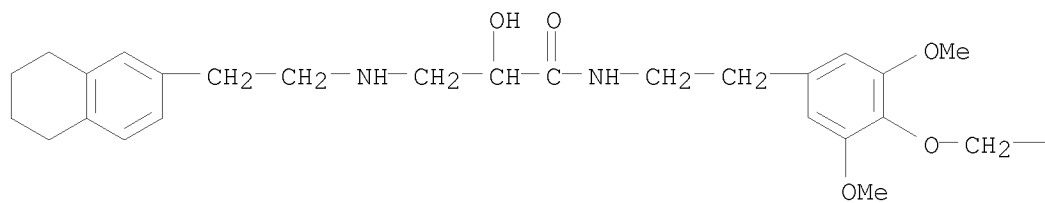


PAGE 1-B

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PAGE 1-A



PAGE 1-B

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10/513699

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L5 ANSWER 6 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2003:1014567 CAPLUS

DOCUMENT NUMBER: 138:411231

TITLE: Preparation of α -oxygenated or α -thiolated carboxylic acid phenethylamides for controlling fungal infestation in plants

INVENTOR(S): Zeller, Martin; Lamberth, Clemens; Kriz, Miroslav

PATENT ASSIGNEE(S): Syngenta Participations A.-G., Switz.

SOURCE: PCT Int. Appl., 100 pp.

CODEN: PIXXD2

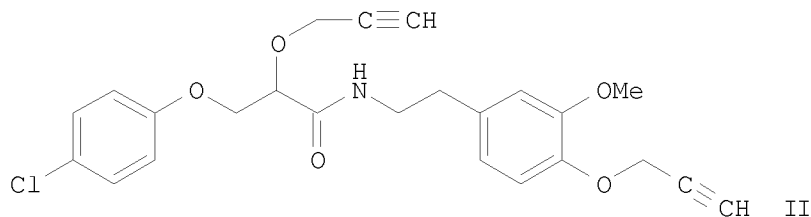
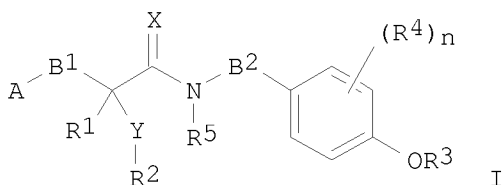
DOCUMENT TYPE: Patent

LANGUAGE: English

PATENT INFORMATION:

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RW:	AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG, TR			
PRIORITY APPLN. INFO.:			GB 2001-27556	20011116

GI



AB Title compds. I [wherein A = (un)substituted (hetero)aryl; X = O or S; Y = O or S; R1 = H, (halo)alkyl, (halo)alkenyl, (halo)alkynyl, or (halo)cycloalkyl; R2 and R3 = independently H, (halo)alkyl, (halo)alkenyl, (halo)alkynyl, (halo)cycloalkyl(alkyl), (halo)alkoxyalkyl, (halo)alkoxyalkenyl, (halo)alkoxyalkynyl, or (un)substituted aryl(oxy)alkyl, arylalkenyl, or arylalkynyl; or R3 = (un)substituted

heteroarylalkyl, heteroarylalkenyl, or heteroarylalkynyl; R4 = halo, CN, NO2, NH2, CHO, CO2H, (halo)alkyl, (halo)alkenyl(oxy), (halo)alkynyl(oxy), (halo)alkoxy(alkyl), (halo)alkylthio, (halo)alkanoyl, (halo)(di)alkylamino, or (halo)alkoxycarbonyl; R5 = H, alkyl, alkenyl, or alkynyl; n = 0-4; B1 = (CR10R11)q or (CHR10R11)rZ(CR12R13)s; q = 2-4; r = 0-3; s = 1-3; r + s = 1-3; Z = O, S, SO, SO2, NR6, CO, OCO, CO2, NR6CO, or CONR6; R6 = H or alkyl; R10-R13 = independently H or alkyl; B2 = alkylene bridge; and optical isomers and mixts. thereof] were prepared These compds. possess useful plant protecting properties and may be employed advantageously in agricultural practice for controlling or preventing the infestation of plants by phytopathogenic microorganisms, especially fungi. For example, 3-chlorolactic acid was coupled with 4-chlorophenol in 3.3N NaOH to give 3-(4-chlorophenoxy)-2-hydroxypropionic acid. Amidation with 2-[3-methoxy-4-[(prop-2-ynyl)oxy]phenyl]ethylamine•HCl in the presence of N,N-diisopropylethylamine in DMF, followed by etherification with propargyl bromide in toluene provided the N-(phenethyl)- α -(propargyloxy)propionamide II. The latter showed residual protective action and residual curative action against fungal infestation by Plasmopara viticola on vines, Phytophthora on tomato plants, and Phytophthora on potato plants by 80-100% at 200 ppm. [This abstract record is one of 6 records for this document necessitated by the large number of index entries required to fully index the document and publication system constraints.]

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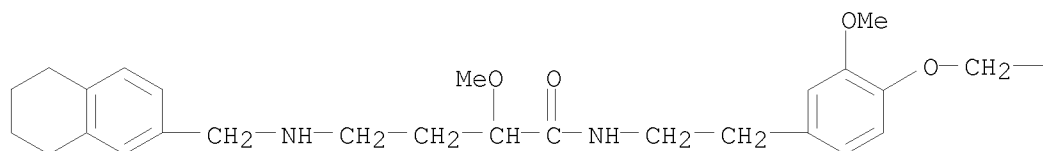
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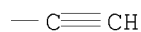
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PAGE 1-A



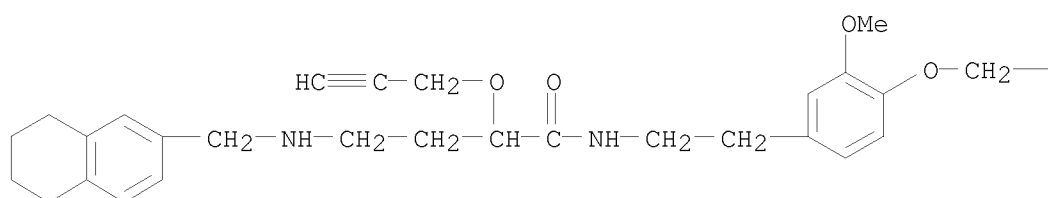
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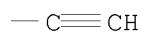
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PAGE 1-A



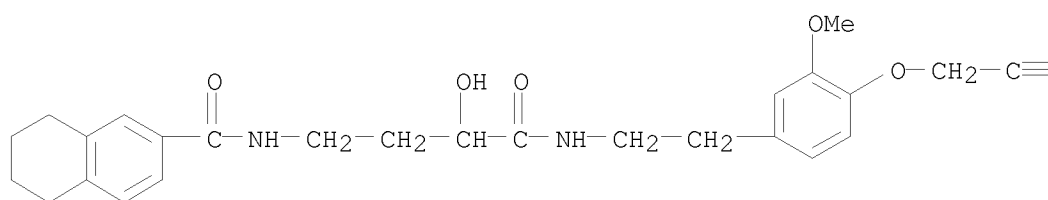
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PAGE 1-A



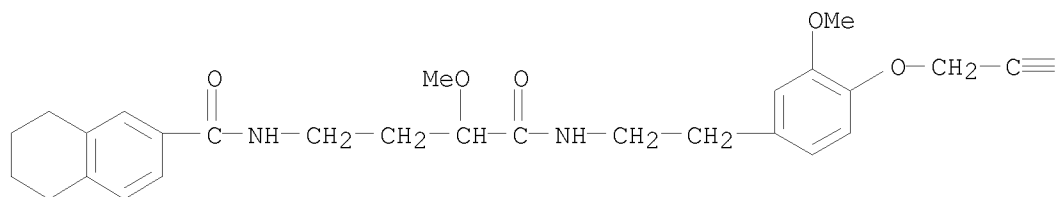
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PAGE 1-A

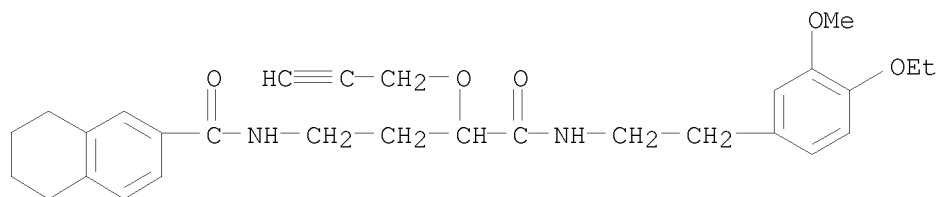


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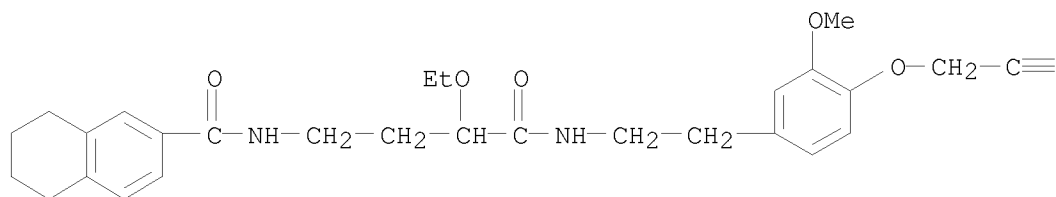
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PAGE 1-A



PAGE 1-B

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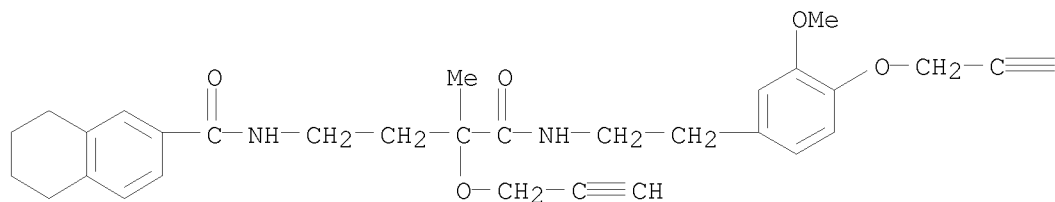
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10/513699

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PAGE 1-A



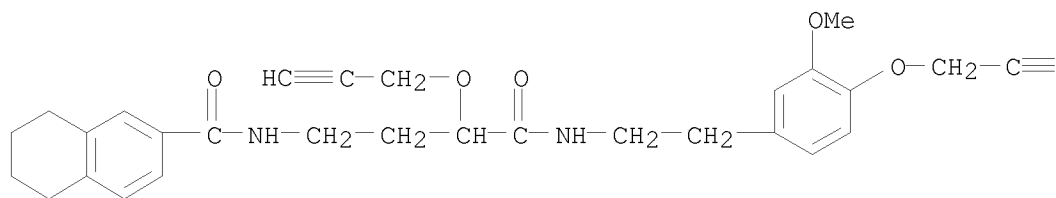
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PAGE 1-A



PAGE 1-B

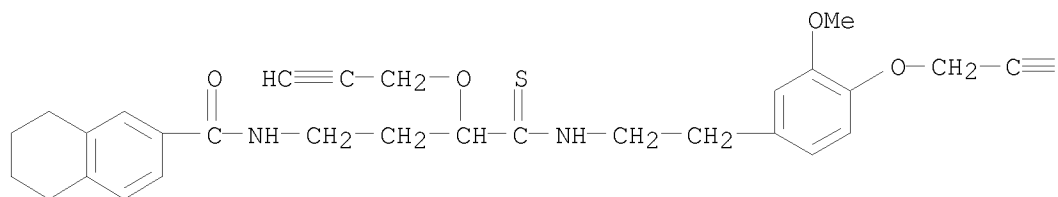
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10/513699

PAGE 1-A



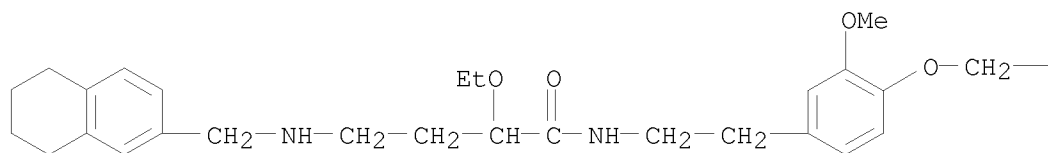
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PAGE 1-A



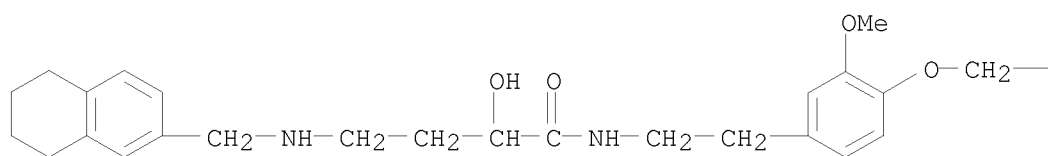
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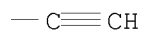
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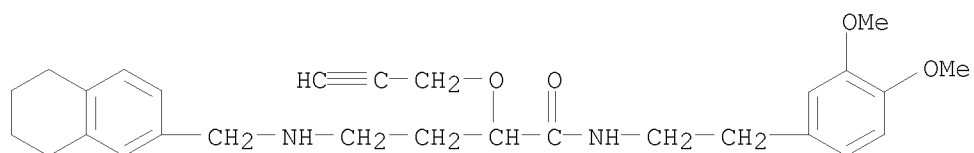


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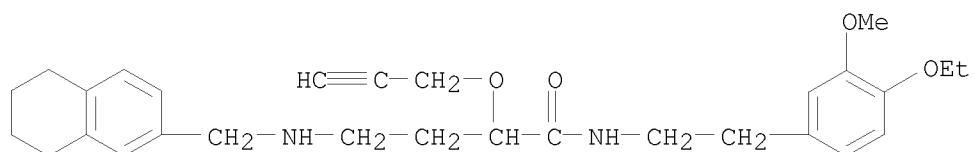


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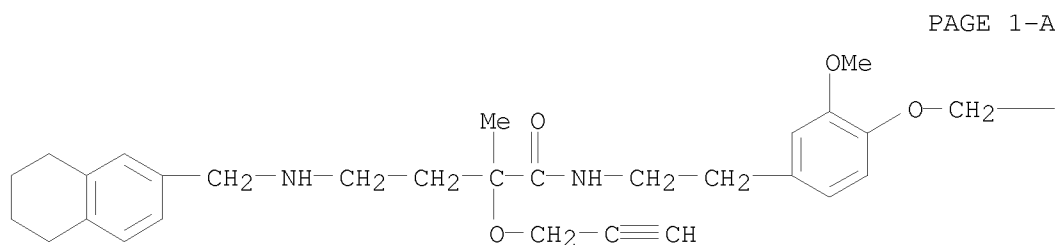
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RN 1067816-90-3 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



RN 1067816-91-4 CAPLUS

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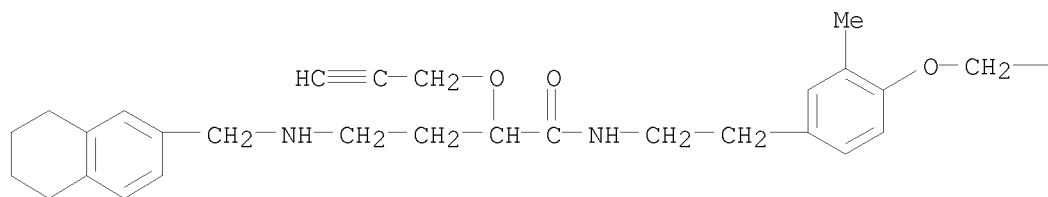
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PAGE 1-A



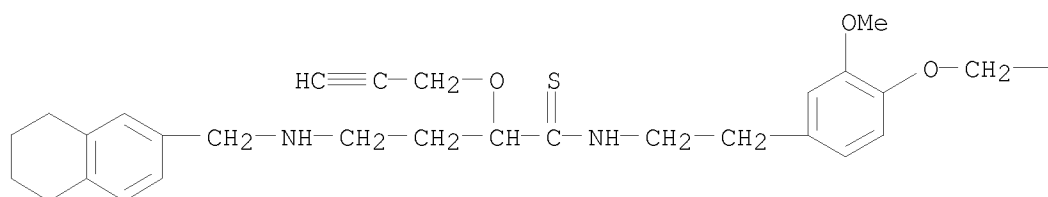
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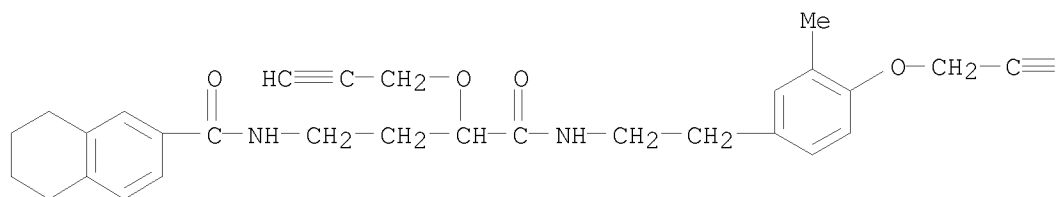


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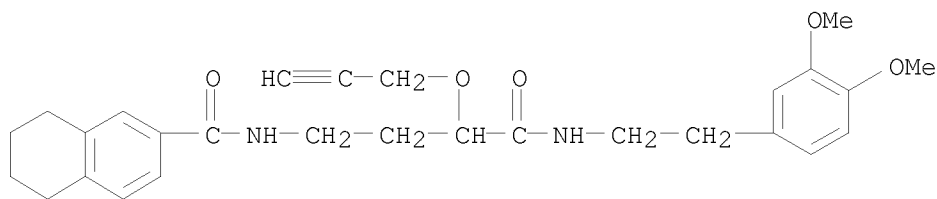
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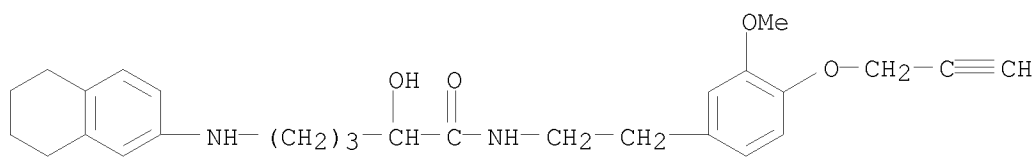
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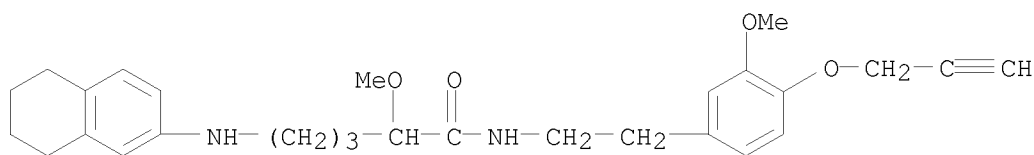
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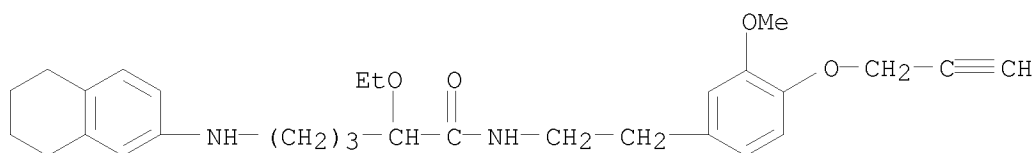
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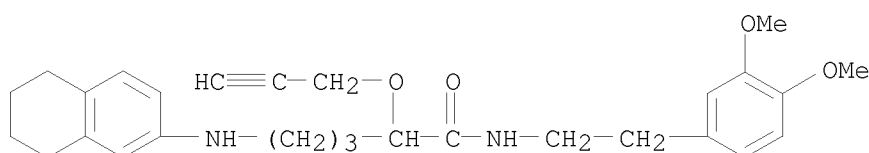
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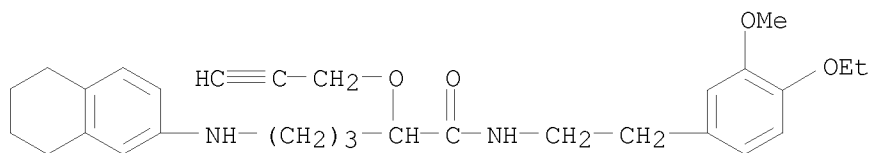
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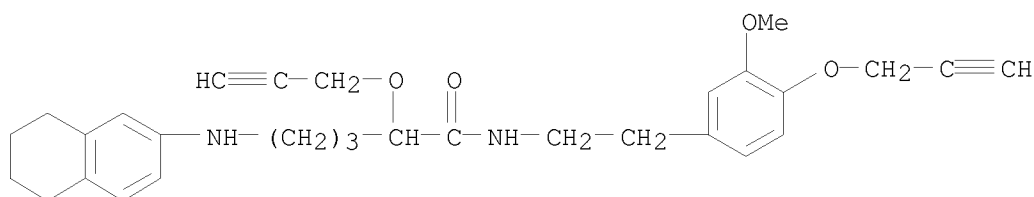
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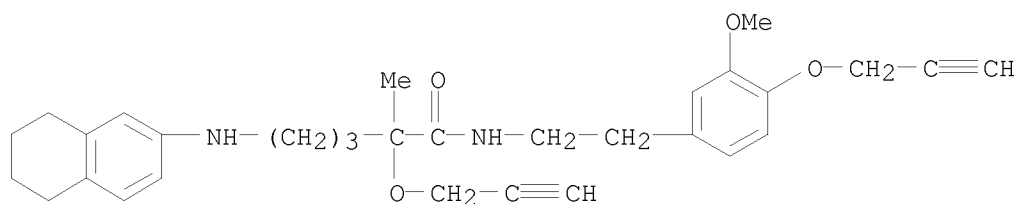
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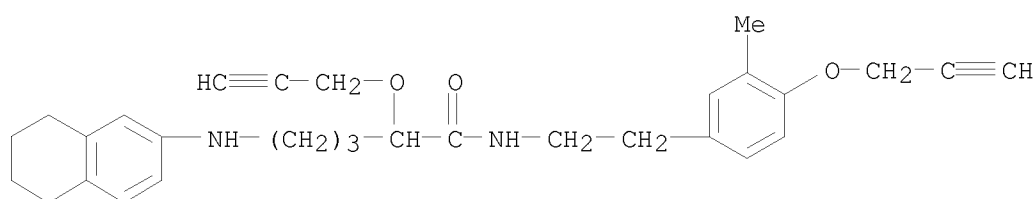
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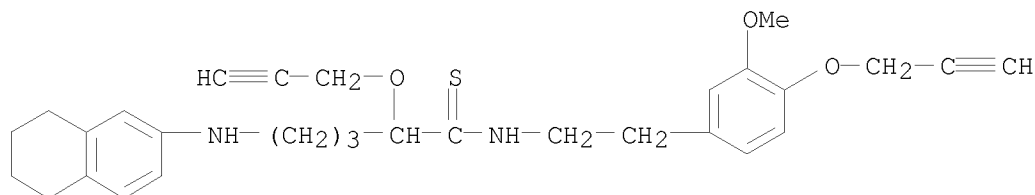
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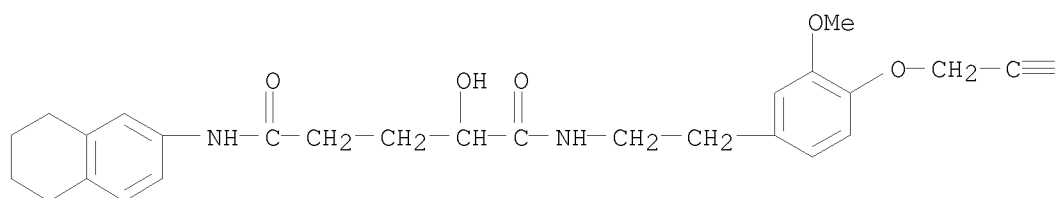
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PAGE 1-A



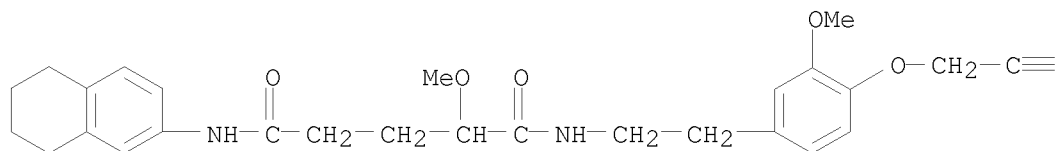
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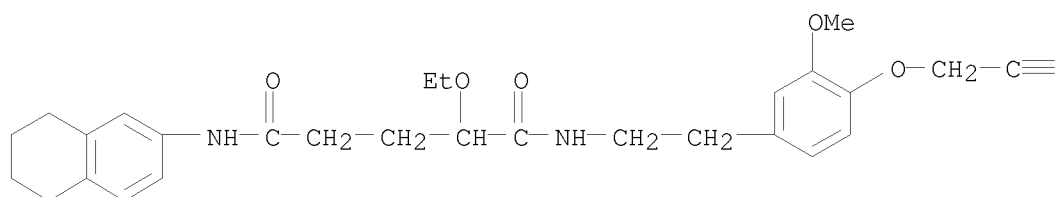
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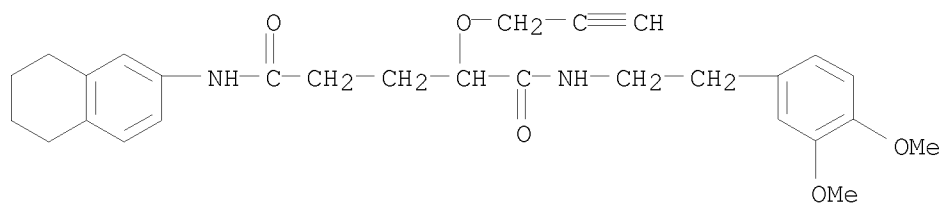
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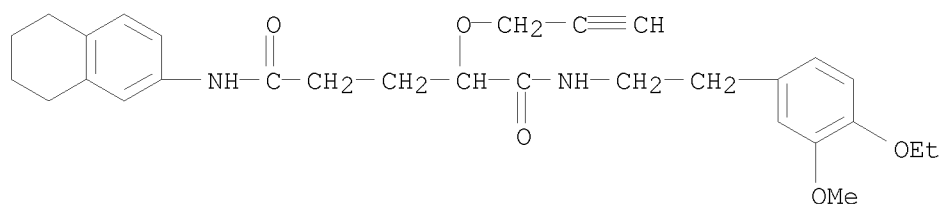
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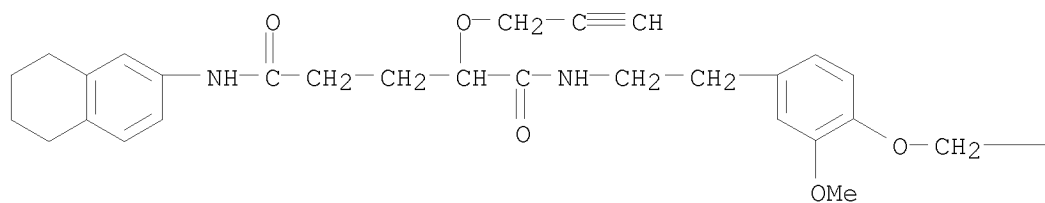


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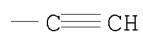


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PAGE 1-A



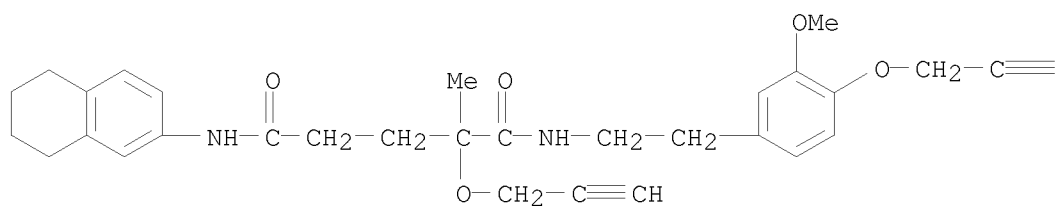
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PAGE 1-A



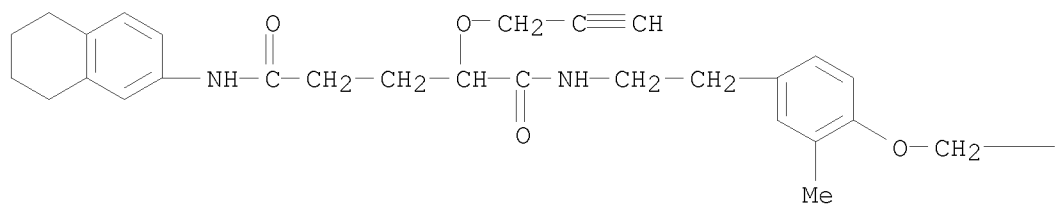
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PAGE 1-A



PAGE 1-B

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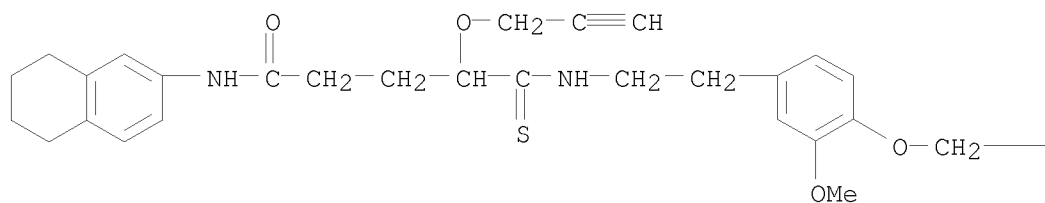
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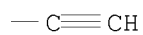
Erich Leese

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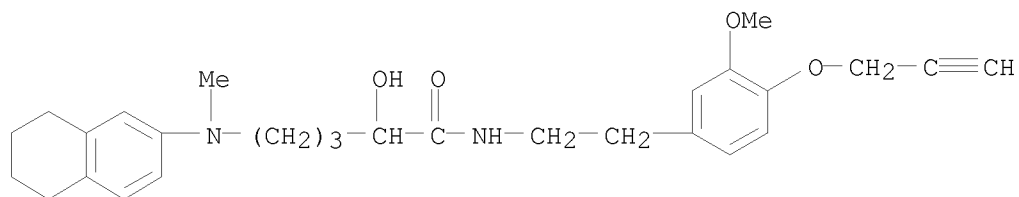
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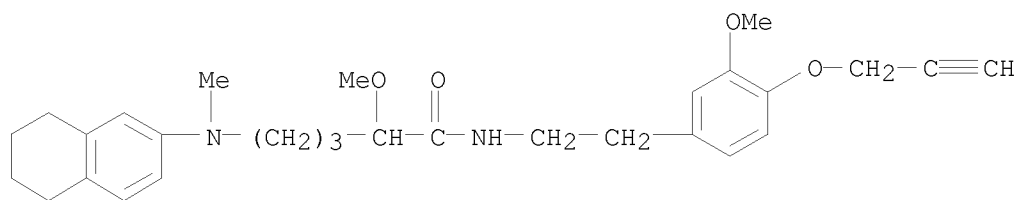
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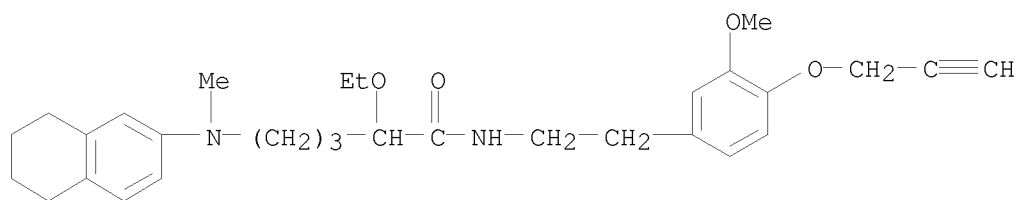


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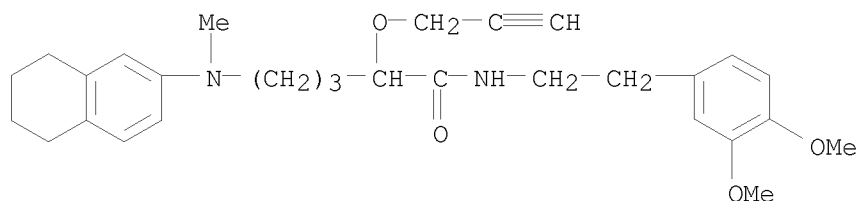
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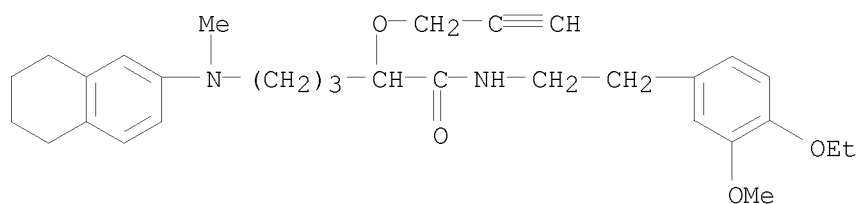
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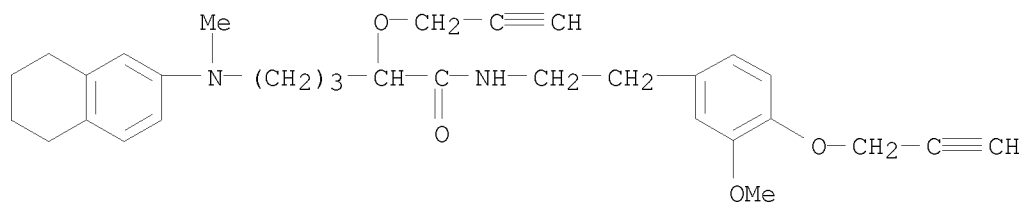
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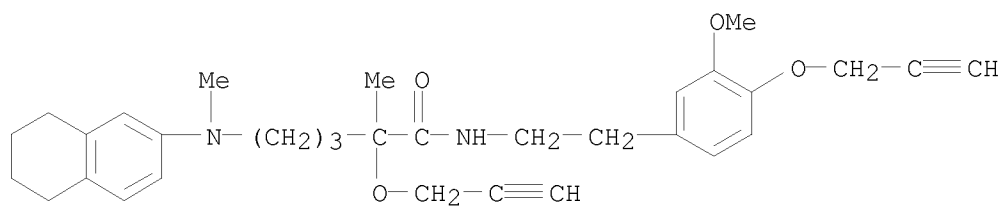


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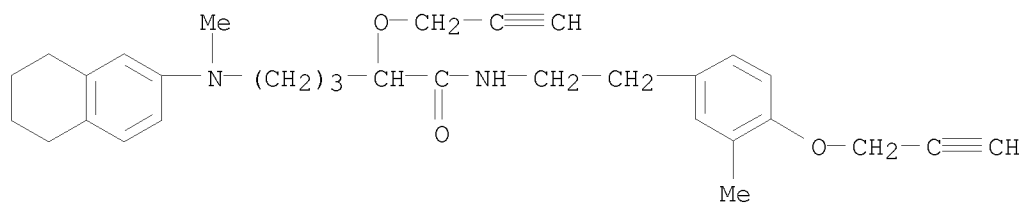


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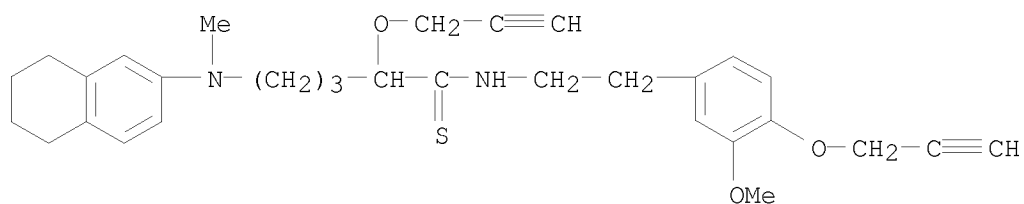
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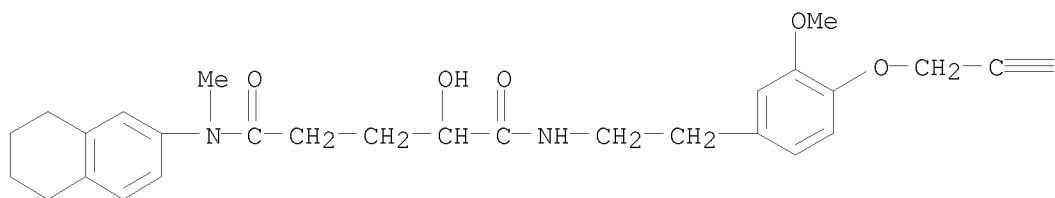


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PAGE 1-A



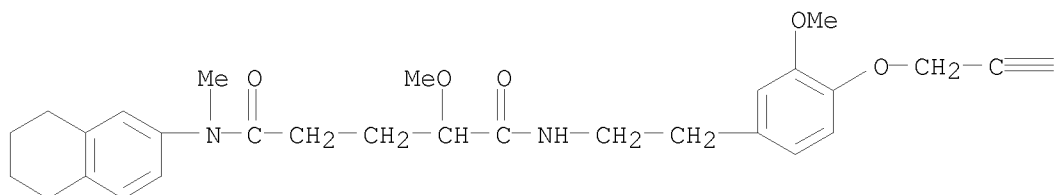
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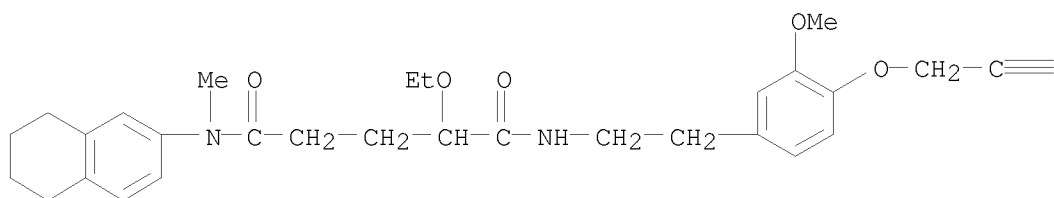
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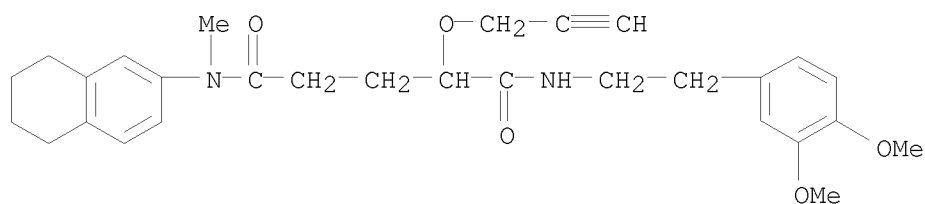
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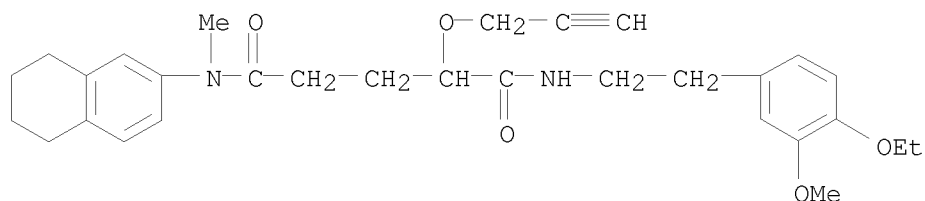
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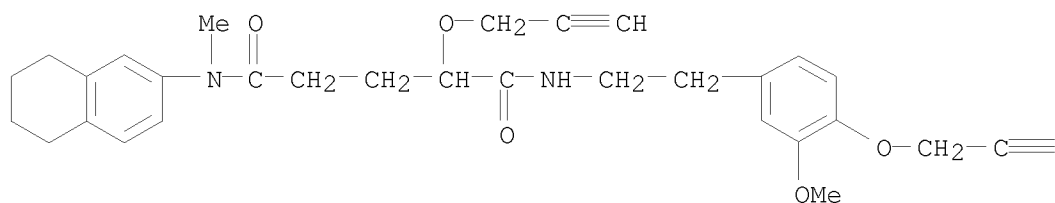


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PAGE 1-A



PAGE 1-B

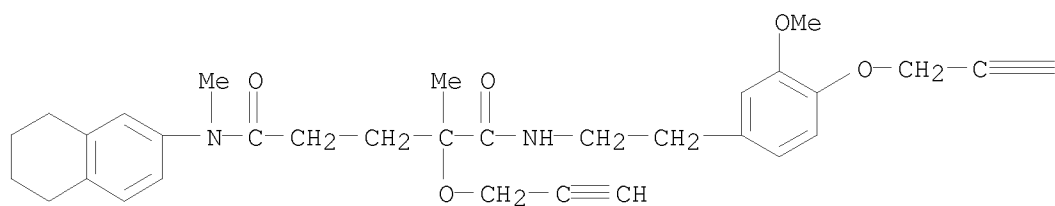
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<12/04/2007>

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PAGE 1-A



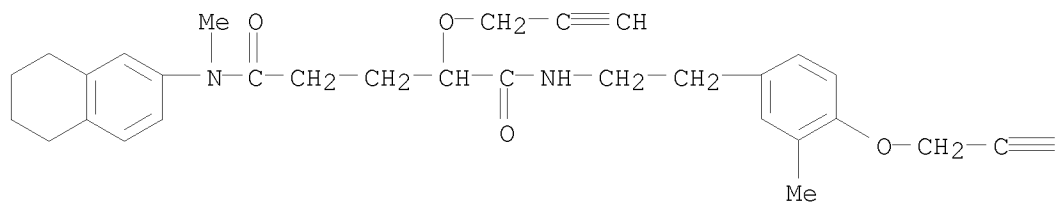
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PAGE 1-B

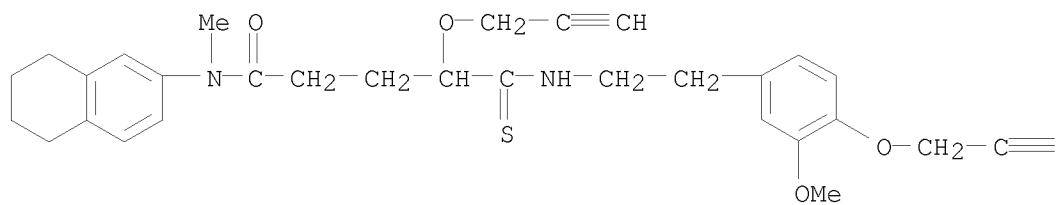
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PAGE 1-A



PAGE 1-B

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REFERENCE COUNT:

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THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS
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ACCESSION NUMBER: 2003:1014566 CAPLUS

DOCUMENT NUMBER: 138:411230

TITLE: Preparation of α -oxygenated or α -thiolated carboxylic acid phenethylamides for controlling fungal infestation in plants

INVENTOR(S): Zeller, Martin; Lamberth, Clemens; Kriz, Miroslav

PATENT ASSIGNEE(S): Syngenta Participations A.-G., Switz.

SOURCE: PCT Int. Appl., 100 pp.

CODEN: PIXXD2

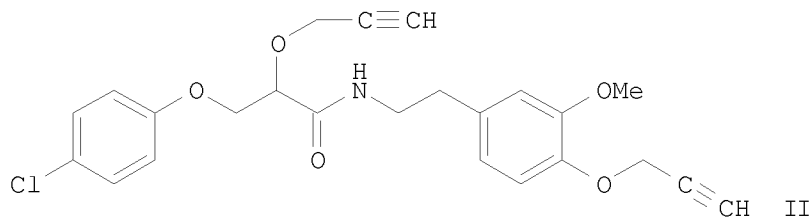
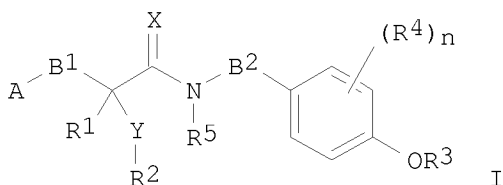
DOCUMENT TYPE: Patent

LANGUAGE: English

PATENT INFORMATION:

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RW:	AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG, TR			
PRIORITY APPLN. INFO.:			GB 2001-27556	20011116

GI



AB Title compds. I [wherein A = (un)substituted (hetero)aryl; X = O or S; Y = O or S; R1 = H, (halo)alkyl, (halo)alkenyl, (halo)alkynyl, or (halo)cycloalkyl; R2 and R3 = independently H, (halo)alkyl, (halo)alkenyl, (halo)alkynyl, (halo)cycloalkyl(alkyl), (halo)alkoxyalkyl, (halo)alkoxyalkenyl, (halo)alkoxyalkynyl, or (un)substituted aryl(oxy)alkyl, arylalkenyl, or arylalkynyl; or R3 = (un)substituted

heteroarylalkyl, heteroarylalkenyl, or heteroarylalkynyl; R4 = halo, CN, NO2, NH2, CHO, CO2H, (halo)alkyl, (halo)alkenyl(oxy), (halo)alkynyl(oxy), (halo)alkoxy(alkyl), (halo)alkylthio, (halo)alkanoyl, (halo)(di)alkylamino, or (halo)alkoxycarbonyl; R5 = H, alkyl, alkenyl, or alkynyl; n = 0-4; B1 = (CR10R11)q or (CHR10R11)rZ(CR12R13)s; q = 2-4; r = 0-3; s = 1-3; r + s = 1-3; Z = O, S, SO, SO2, NR6, CO, OCO, CO2, NR6CO, or CONR6; R6 = H or alkyl; R10-R13 = independently H or alkyl; B2 = alkylene bridge; and optical isomers and mixts. thereof] were prepared These compds. possess useful plant protecting properties and may be employed advantageously in agricultural practice for controlling or preventing the infestation of plants by phytopathogenic microorganisms, especially fungi. For example, 3-chlorolactic acid was coupled with 4-chlorophenol in 3.3N NaOH to give 3-(4-chlorophenoxy)-2-hydroxypropionic acid. Amidation with 2-[3-methoxy-4-[(prop-2-ynyl)oxy]phenyl]ethylamine•HCl in the presence of N,N-diisopropylethylamine in DMF, followed by etherification with propargyl bromide in toluene provided the N-(phenethyl)- α -(propargyloxy)propionamide II. The latter showed residual protective action and residual curative action against fungal infestation by Plasmopara viticola on vines, Phytophthora on tomato plants, and Phytophthora on potato plants by 80-100% at 200 ppm. [This abstract record is one of 6 records for this document necessitated by the large number of index entries required to fully index the document and publication system constraints.]

IT	1067726-14-0	1067726-15-1	1067726-16-2
	1067726-17-3	1067726-18-4	1067726-19-5
	1067726-20-8	1067726-21-9	1067726-22-0
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	1067735-19-6	1067735-20-9	1067735-21-0
	1067735-22-1	1067735-23-2	1067735-24-3
	1067751-02-3	1067751-04-5	1067751-05-6
	1067751-06-7	1067751-07-8	1067751-08-9
	1067751-09-0	1067751-10-3	1067751-11-4
	1067751-39-6	1067751-40-9	1067751-41-0
	1067751-42-1	1067751-43-2	1067751-44-3
	1067751-45-4	1067751-46-5	1067751-47-6
	1067753-93-8	1067753-94-9	1067753-95-0
	1067753-96-1	1067753-97-2	1067753-98-3
	1067753-99-4	1067754-00-0	1067754-01-1
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	1067769-34-9	1067770-20-0	1067770-22-2
	1067770-23-3	1067771-37-2	1067771-39-4
	1067772-27-3	1067772-28-4	1067773-43-6
	1067773-44-7	1067773-45-8	1067774-60-0
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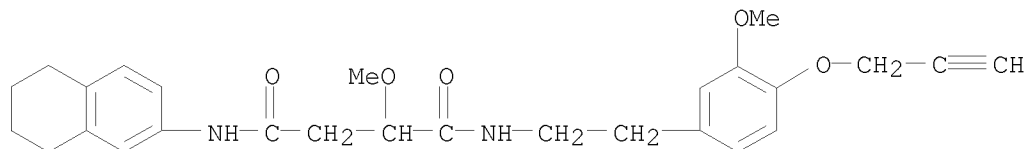
RL: PRPH (Prophetic)

(Preparation of α -oxygenated or α -thiolated carboxylic acid phenethylamides for controlling fungal infestation in plants)

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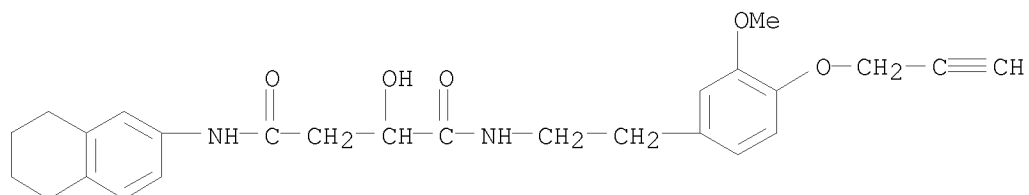
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10/513699



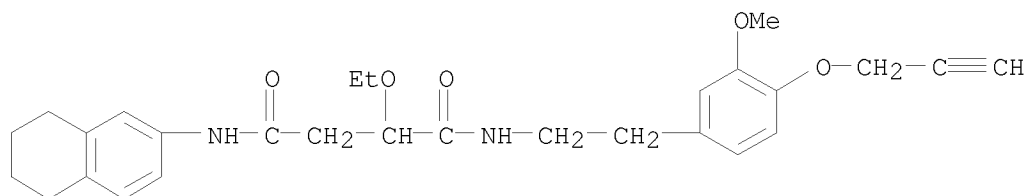
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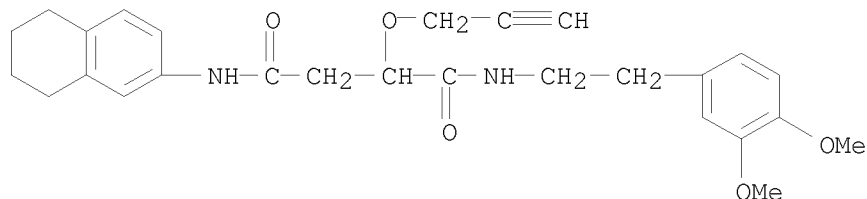
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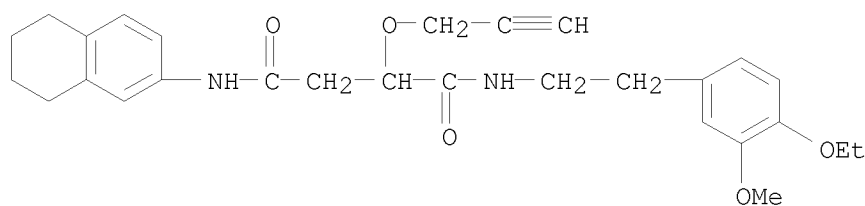
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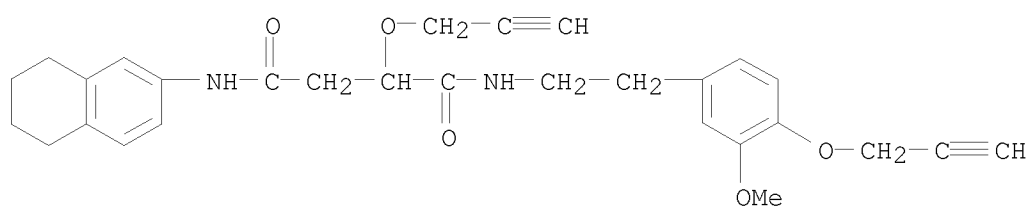
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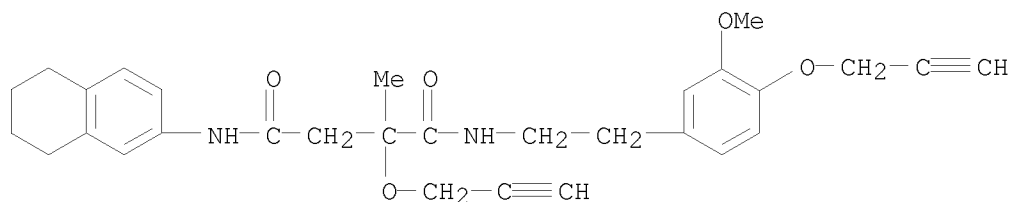
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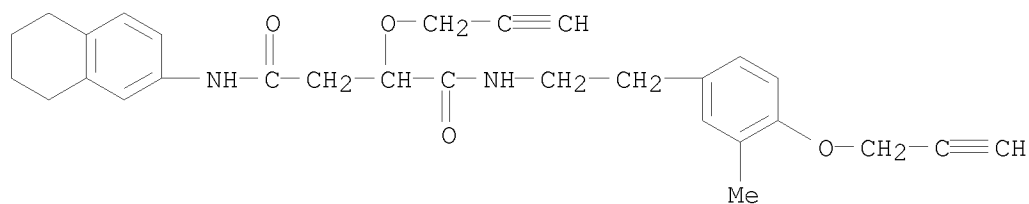
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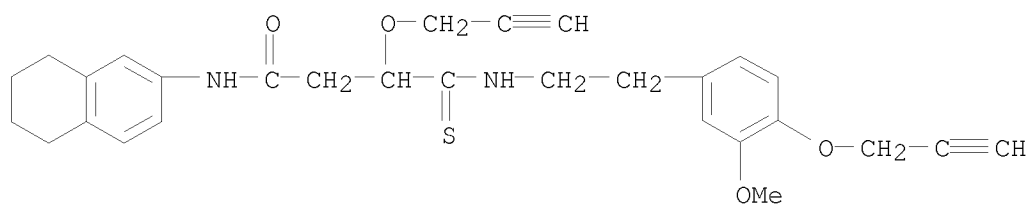
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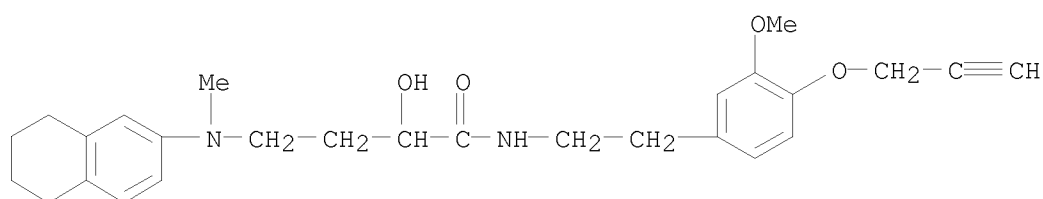
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Erich Leese

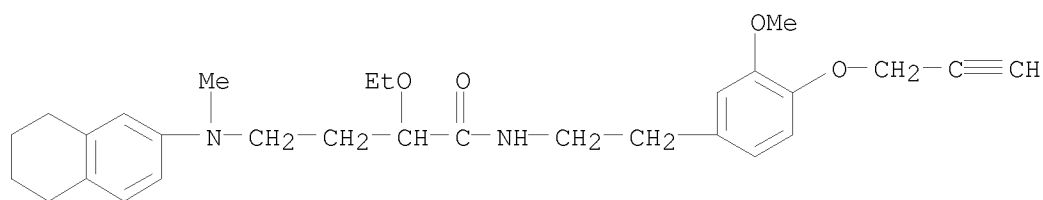
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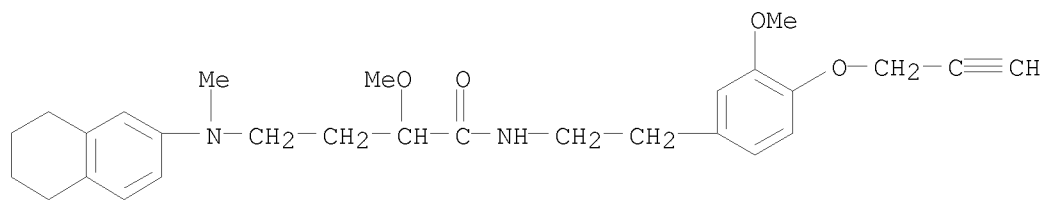
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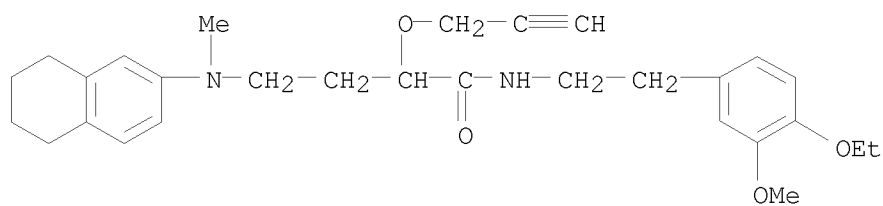


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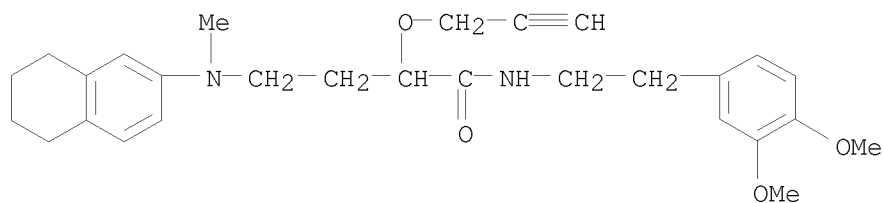


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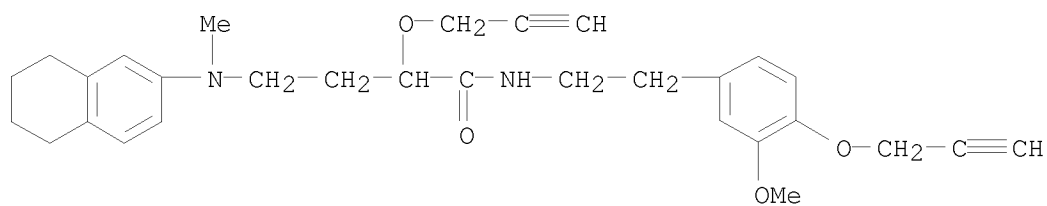
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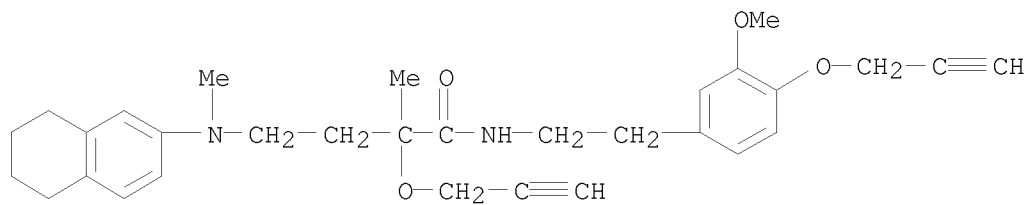
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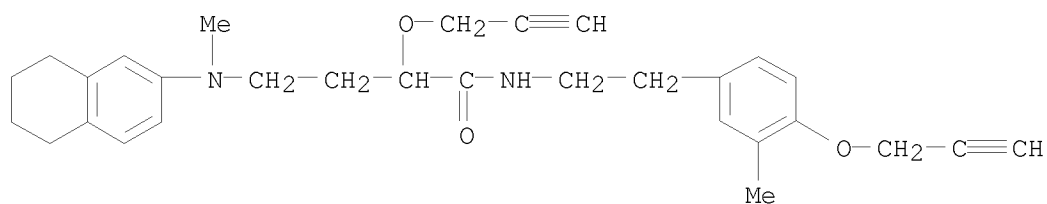


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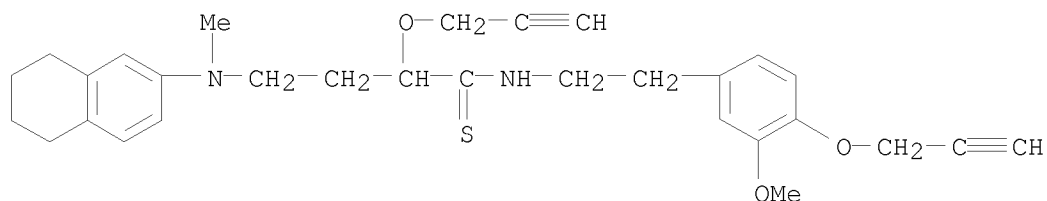


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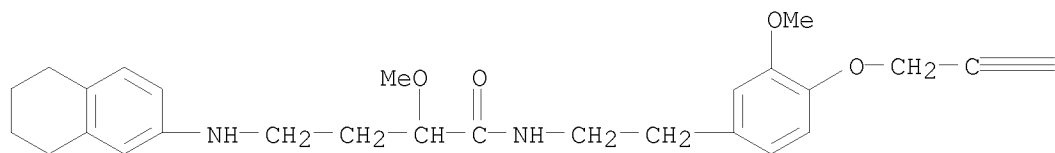


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PAGE 1-A



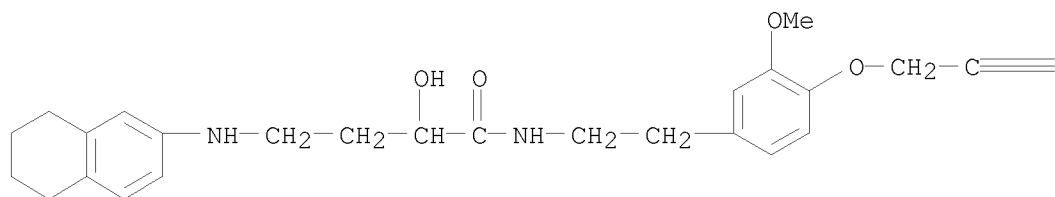
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10/513699

PAGE 1-A

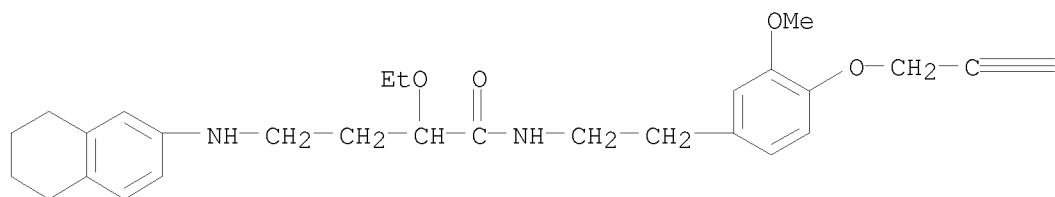


PAGE 1-B

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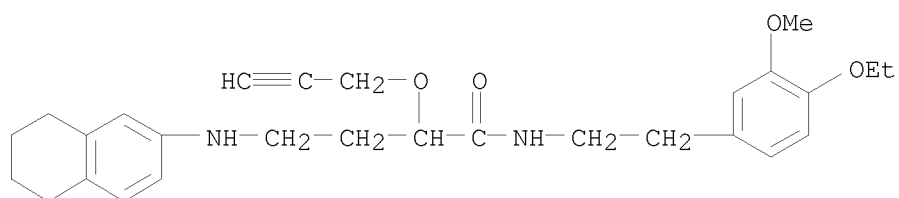
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PAGE 1-B

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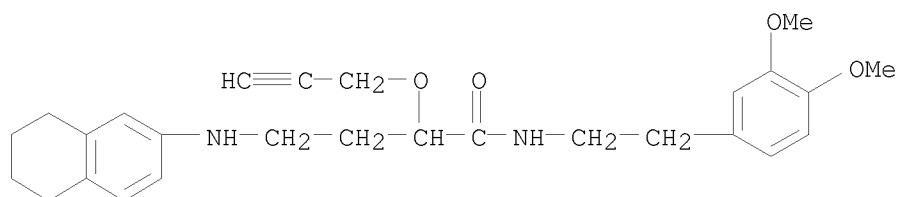


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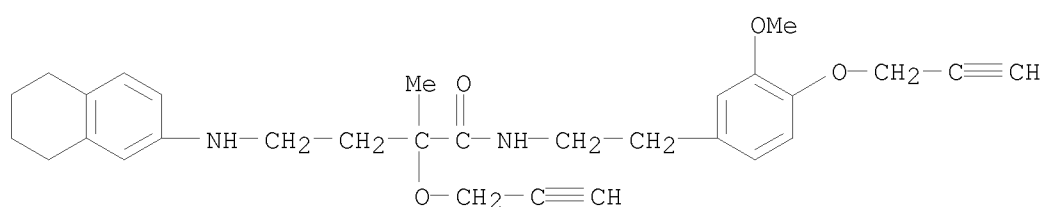
Erich Leese

10/513699



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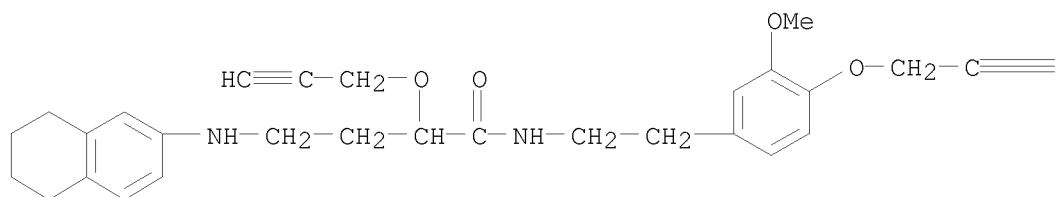
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PAGE 1-A



PAGE 1-B

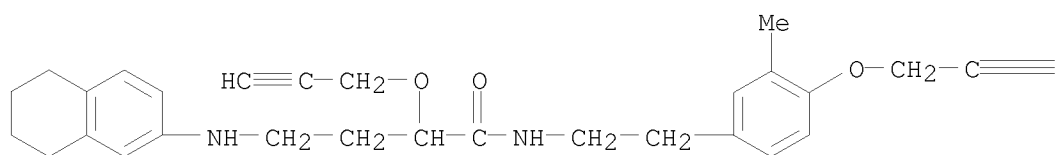
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10/513699

PAGE 1-A



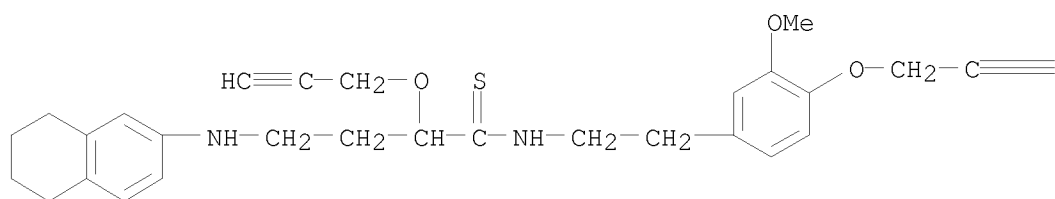
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PAGE 1-A

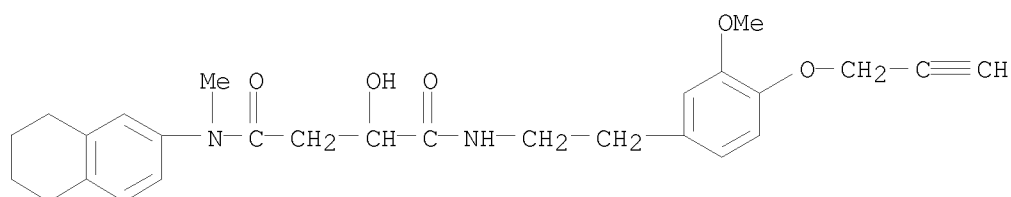


PAGE 1-B

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RN 1067751-39-6 CAPLUS

CN Butanediamide, 2-hydroxy-N1-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-N4-methyl-N4-(5,6,7,8-tetrahydro-2-naphthalenyl)- (CA INDEX NAME)



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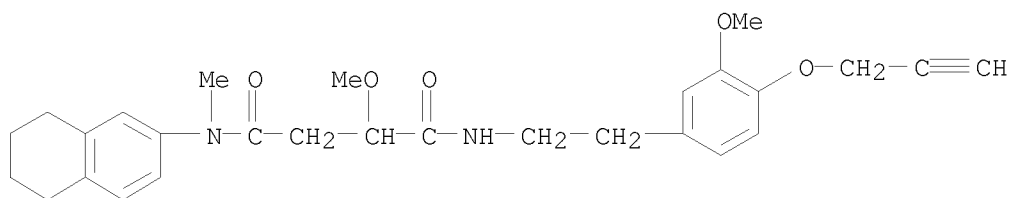
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<12/04/2007>

Erich Leese

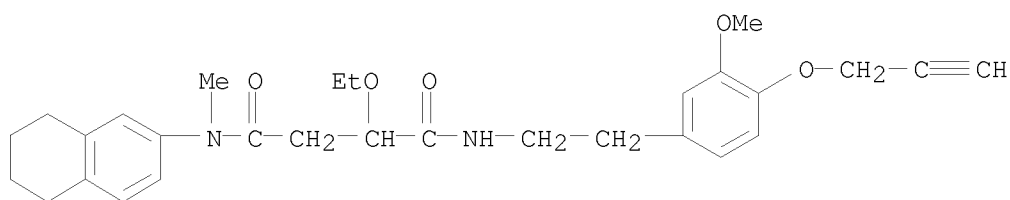
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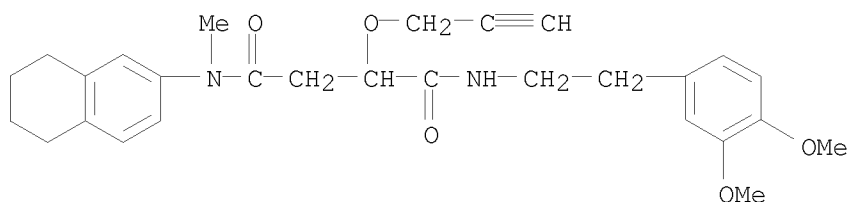
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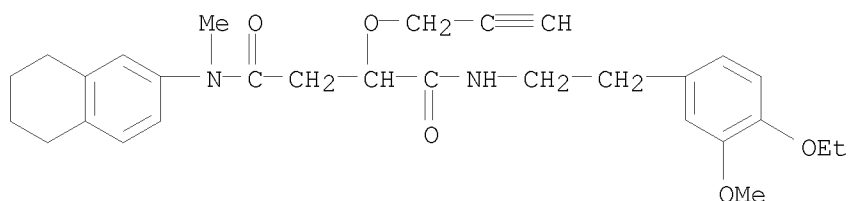
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CN INDEX NAME NOT YET ASSIGNED



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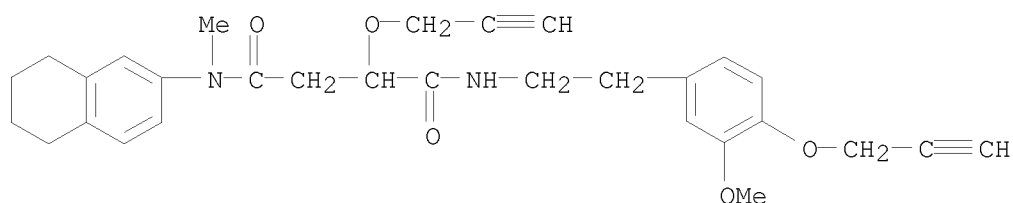
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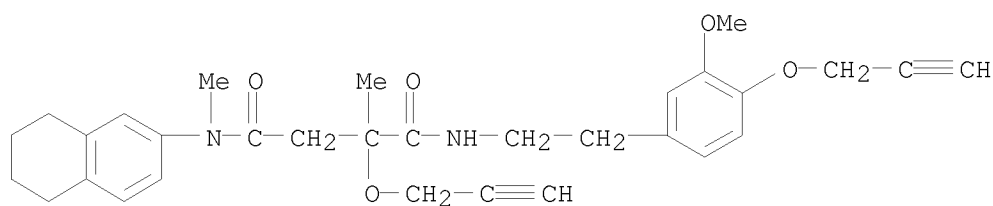
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INDEX NAME)



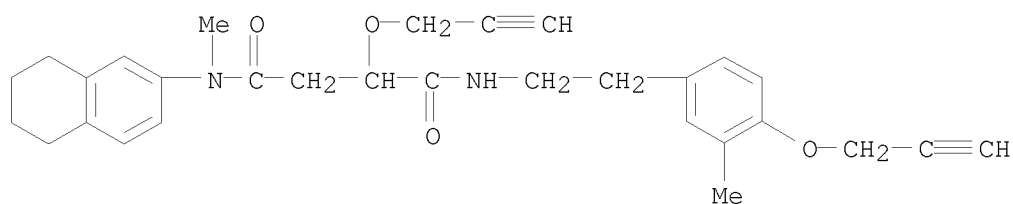
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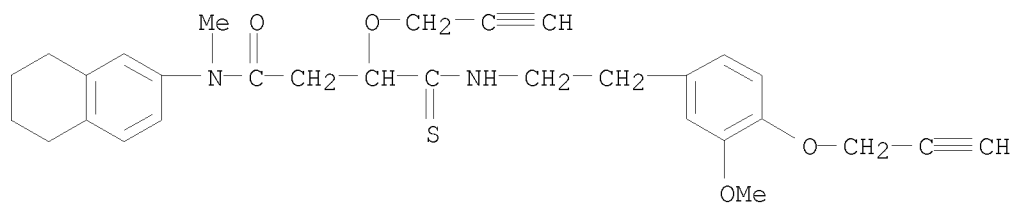
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RN 1067751-47-6 CAPLUS

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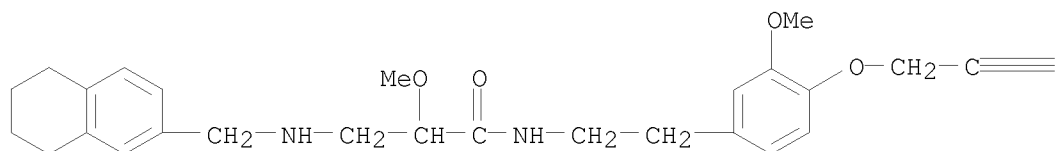
Erich Leese

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RN 1067753-93-8 CAPLUS

CN Propanamide, 2-methoxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-3-
[[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]amino]- (CA INDEX NAME)

PAGE 1-A



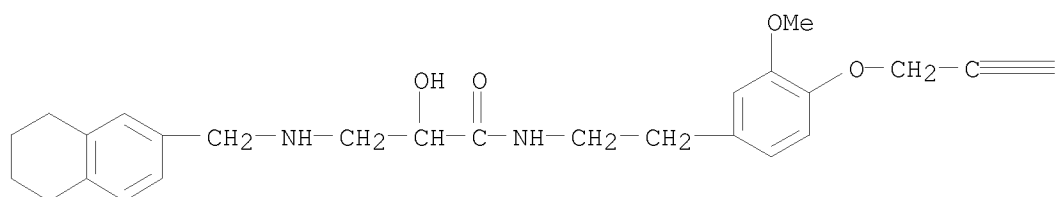
PAGE 1-B

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PAGE 1-A



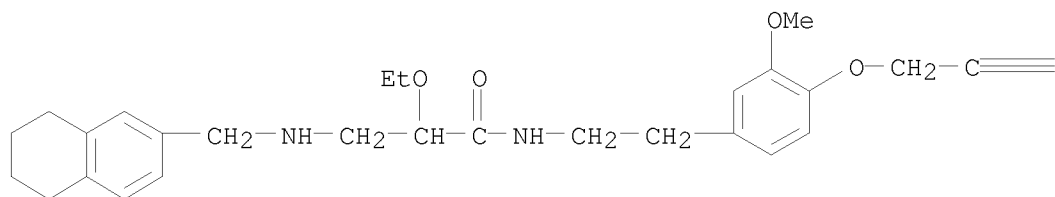
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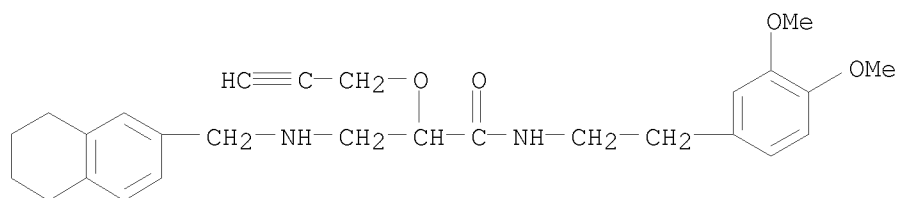
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PAGE 1-A

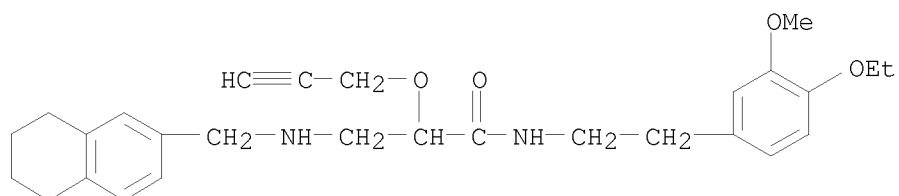


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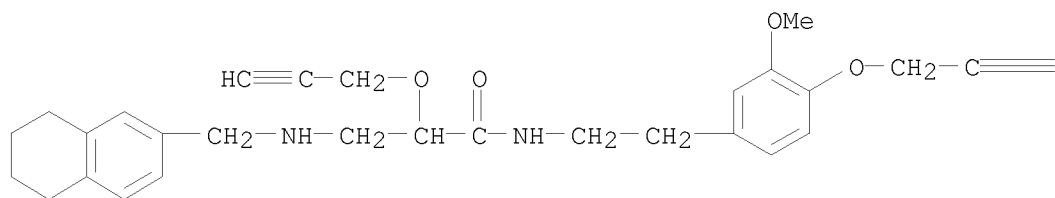
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RN 1067753-97-2 CAPLUS

CN Propanamide, N-[2-(4-ethoxy-3-methoxyphenyl)ethyl]-2-(2-propyn-1-yloxy)-3-
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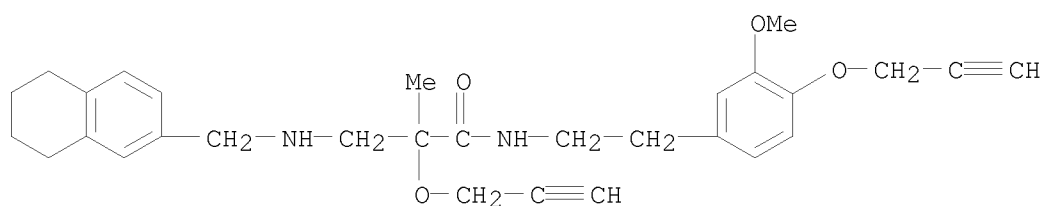
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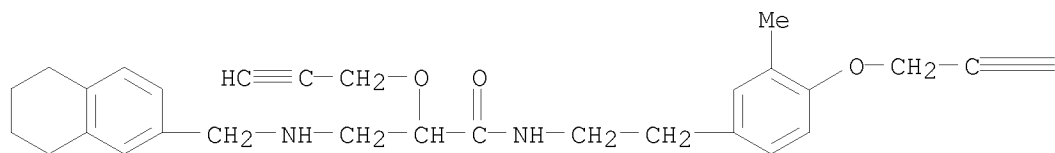
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CN Propanamide, N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-2-methyl-2-(2-propyn-1-yloxy)-3-[[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]amino]-
(CA INDEX NAME)



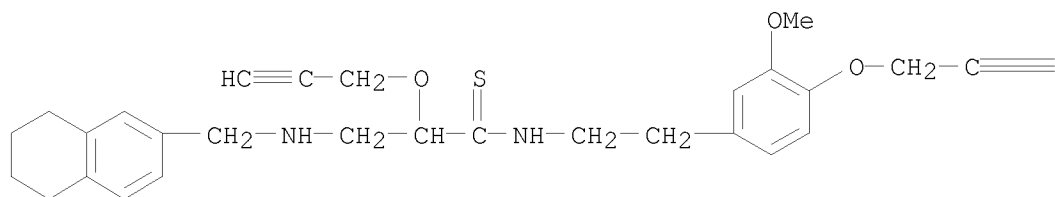
RN 1067754-00-0 CAPLUS

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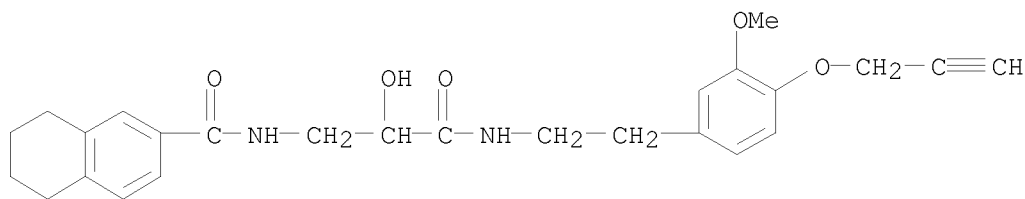
RN 1067754-01-1 CAPLUS

CN Propanethioamide, N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-2-(2-propyn-1-yloxy)-3-[[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]amino]- (CA INDEX NAME)


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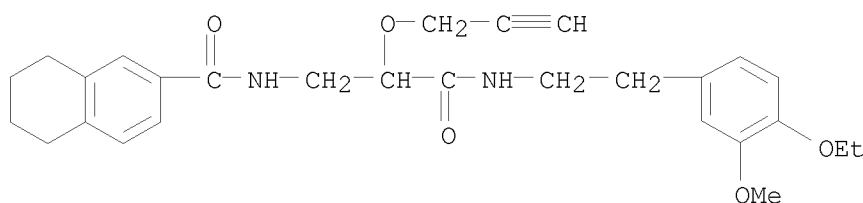
RN 1067766-98-6 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[2-hydroxy-3-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-3-oxopropyl]- (CA INDEX NAME)



RN 1067766-99-7 CAPLUS

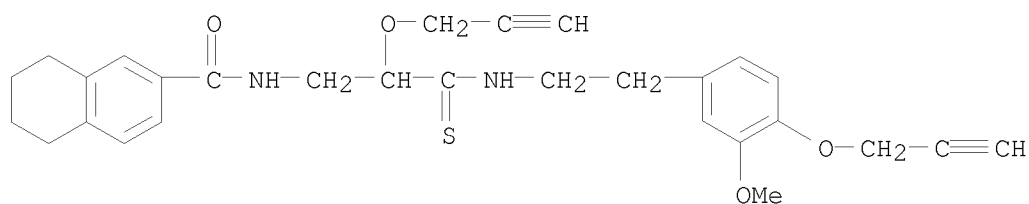
CN 2-Naphthalenecarboxamide, N-[3-[[2-(4-ethoxy-3-methoxyphenyl)ethyl]amino]-3-oxo-2-(2-propyn-1-yloxy)propyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 1067767-00-3 CAPLUS

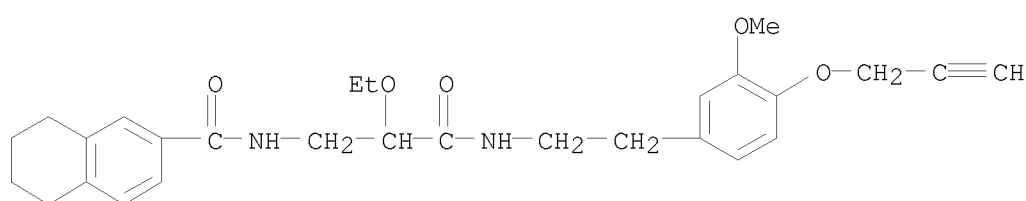
CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[3-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-2-(2-propyn-1-yloxy)-3-thioxopropyl]- (CA INDEX NAME)

10/513699



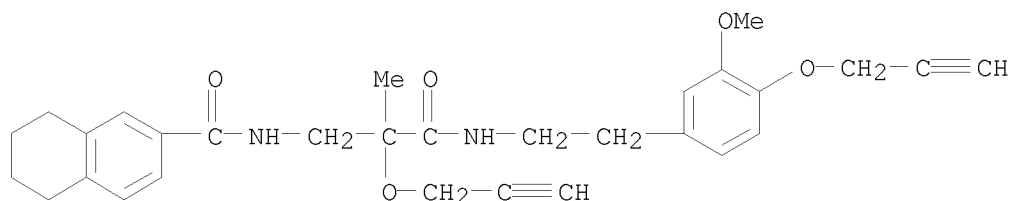
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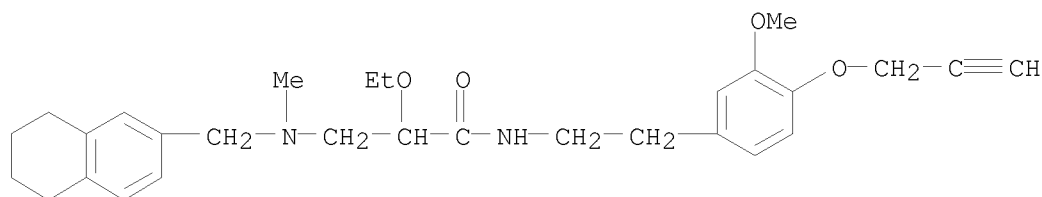
RN 1067767-88-7 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[3-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-2-methyl-3-oxo-2-(2-propyn-1-yloxy)propyl]- (CA INDEX NAME)



RN 1067769-33-8 CAPLUS

CN Propanamide, 2-ethoxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-3-[methyl[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]amino]- (CA INDEX NAME)

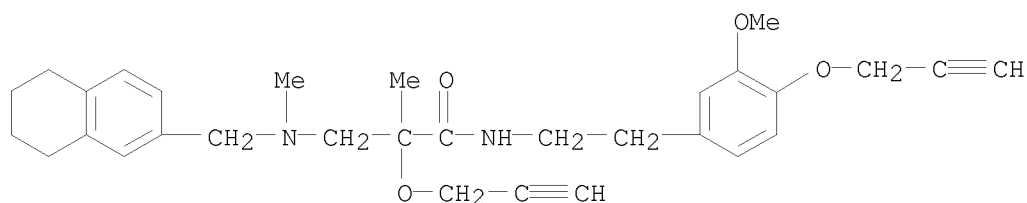


RN 1067769-34-9 CAPLUS

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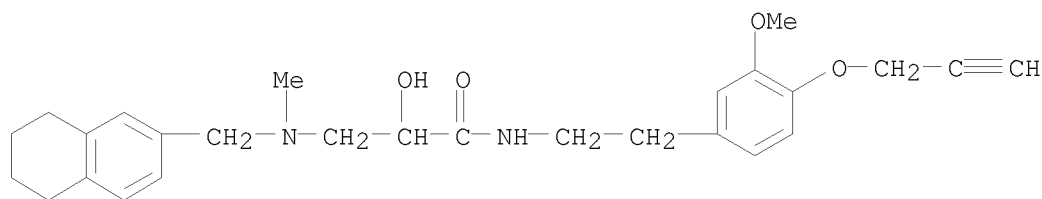
10/513699

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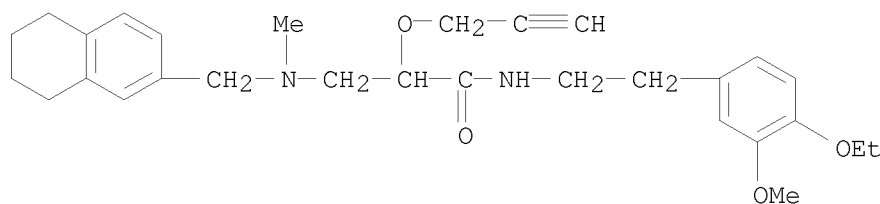
RN 1067770-20-0 CAPLUS

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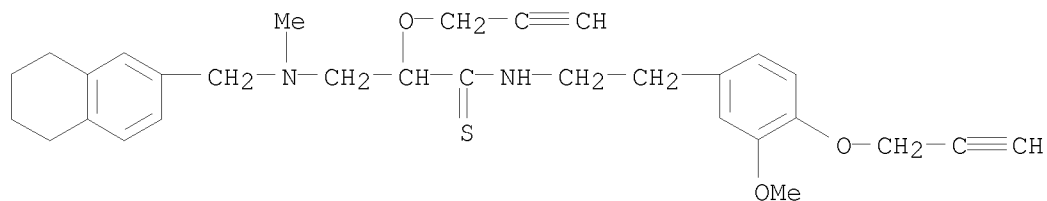
RN 1067770-22-2 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



RN 1067770-23-3 CAPLUS

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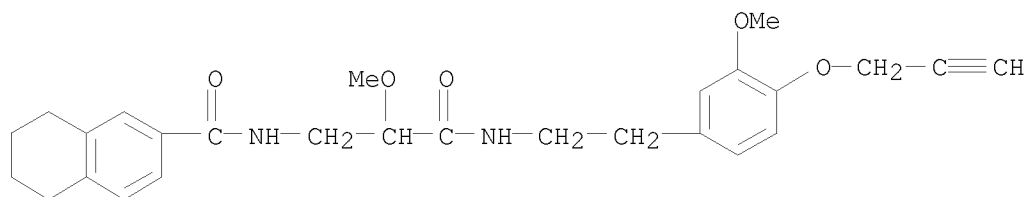
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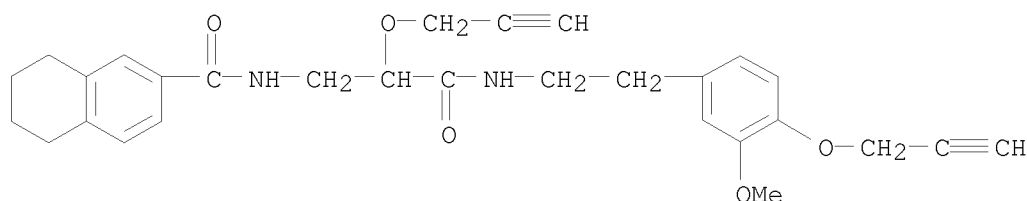
Erich Leese

10/513699



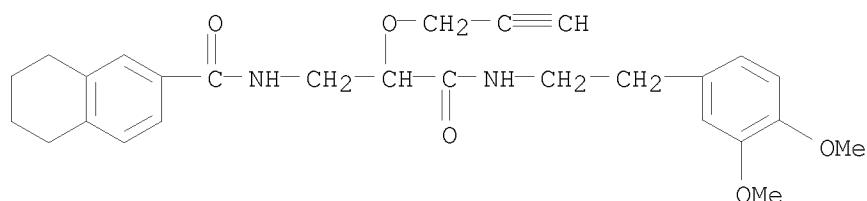
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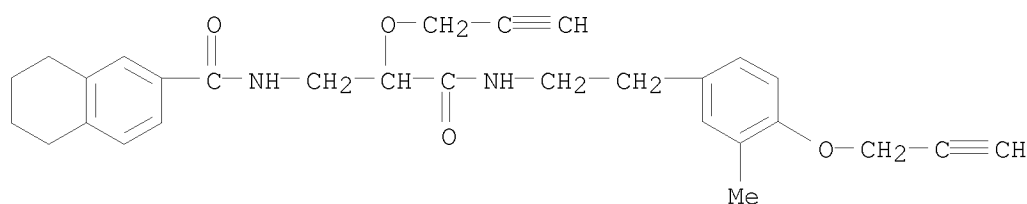
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RN 1067772-28-4 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[3-[[2-[3-methyl-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-3-oxo-2-(2-propyn-1-yloxy)propyl]- (CA INDEX NAME)

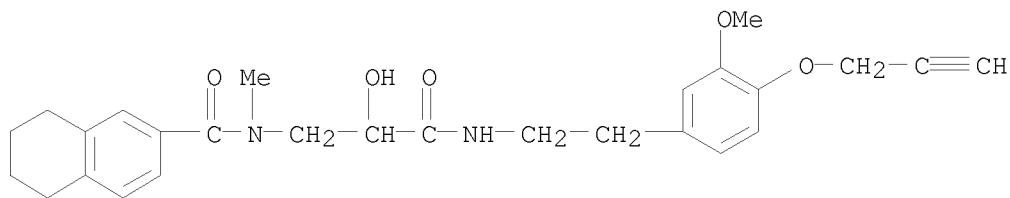


RN 1067773-43-6 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[2-hydroxy-3-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-3-oxo-2-(2-propyn-1-yloxy)propyl]- (CA INDEX NAME)

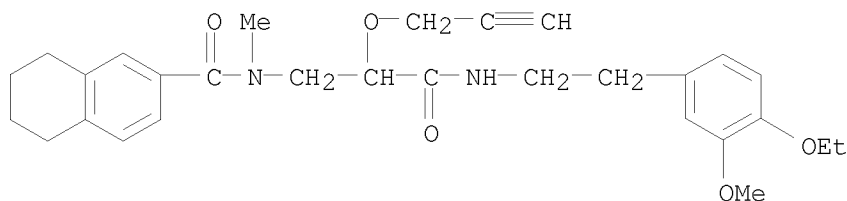
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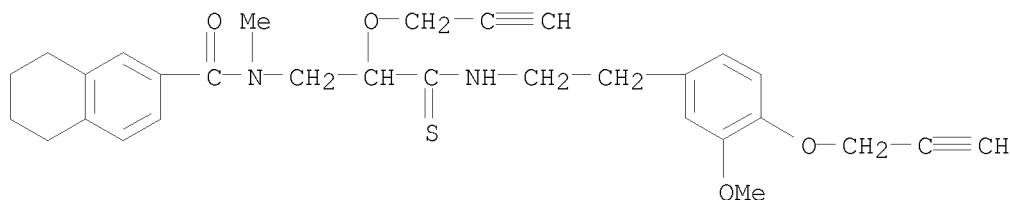
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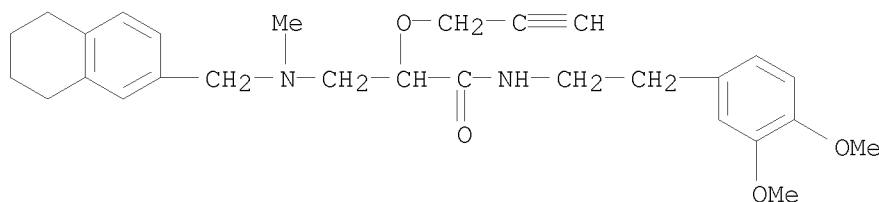
RN 1067773-45-8 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[3-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-2-(2-propyn-1-yloxy)-3-thioxopropyl]-N-methyl- (CA INDEX NAME)



RN 1067774-60-0 CAPLUS

CN Propanamide, N-[2-(3,4-dimethoxyphenyl)ethyl]-3-[methyl[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]amino]-2-(2-propyn-1-yloxy)- (CA INDEX NAME)



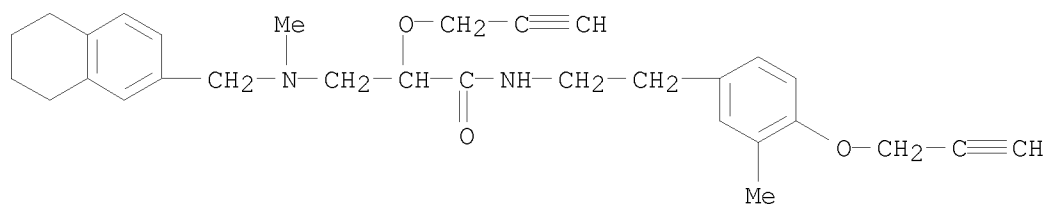
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Erich Leese

10/513699

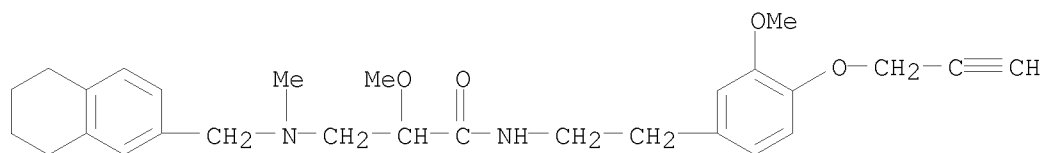
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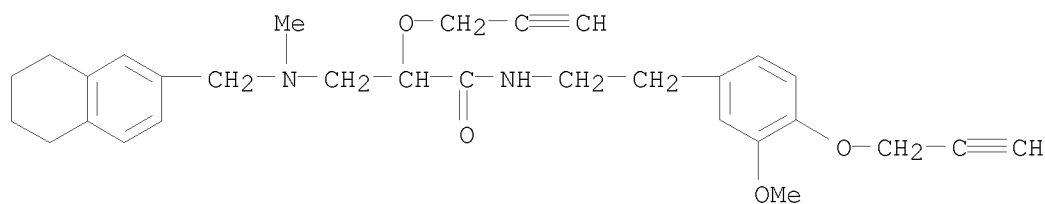
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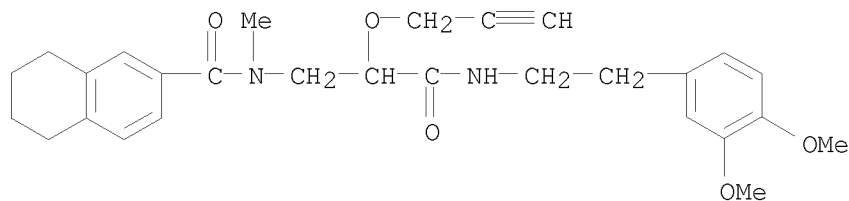
RN 1067775-49-8 CAPLUS

CN Propanamide, N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-3-[methyl[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]amino]-2-(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1067780-47-5 CAPLUS

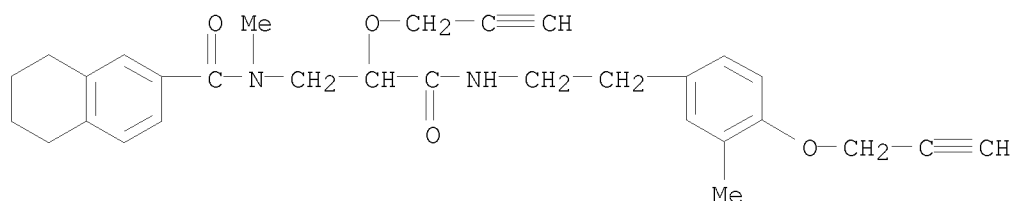
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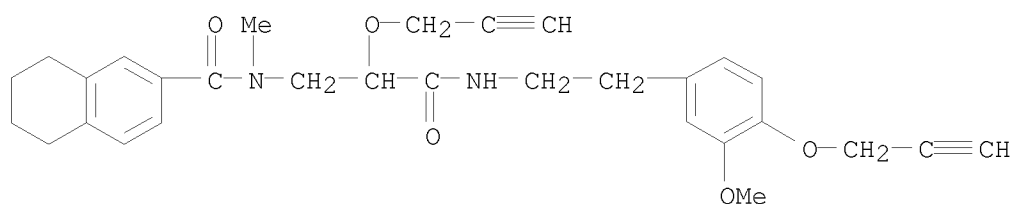
RN 1067780-48-6 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-methyl-N-[3-[[2-[3-methyl-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-3-oxo-2-(2-propyn-1-yloxy)propyl]-
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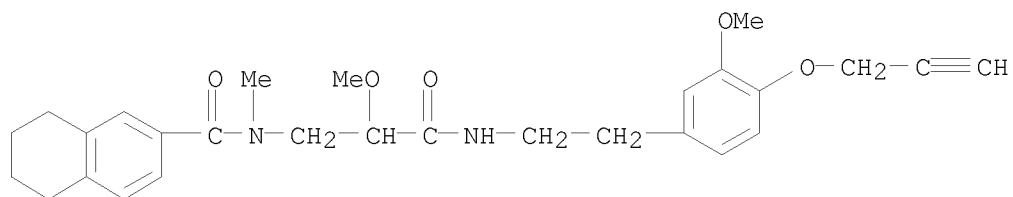
RN 1067782-21-1 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[3-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-3-oxo-2-(2-propyn-1-yloxy)propyl]-N-methyl- (CA INDEX NAME)



RN 1067782-23-3 CAPLUS

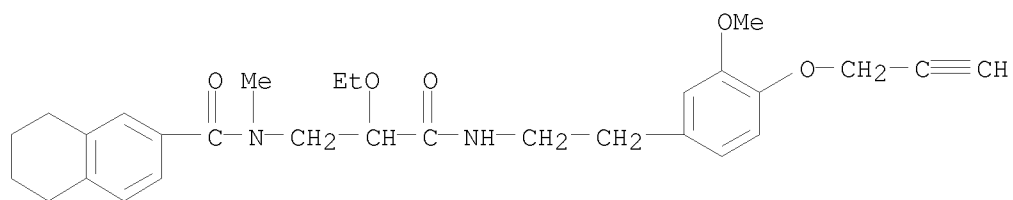
CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[2-methoxy-3-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-3-oxopropyl]-N-methyl- (CA INDEX NAME)



RN 1067784-08-0 CAPLUS

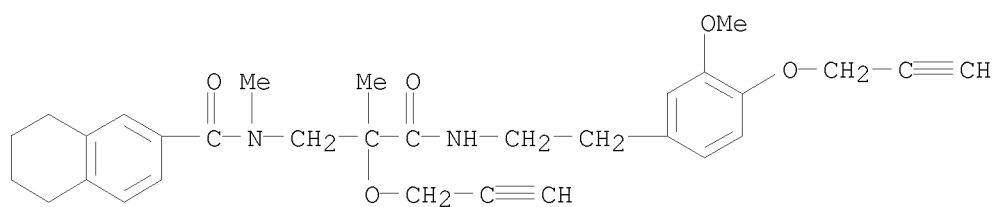
CN 2-Naphthalenecarboxamide, N-[2-ethoxy-3-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-3-oxopropyl]-5,6,7,8-tetrahydro-N-methyl- (CA INDEX NAME)

10/513699



RN 1067784-10-4 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N-[3-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-2-methyl-3-oxo-2-(2-propyn-1-yloxy)propyl]-N-methyl- (CA INDEX NAME)



REFERENCE COUNT:

7

THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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L5 ANSWER 8 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2003:1014565 CAPLUS

DOCUMENT NUMBER: 138:411229

TITLE: Preparation of α -oxygenated or α -thiolated carboxylic acid phenethylamides for controlling fungal infestation in plants

INVENTOR(S): Zeller, Martin; Lamberth, Clemens; Kriz, Miroslav

PATENT ASSIGNEE(S): Syngenta Participations A.-G., Switz.

SOURCE: PCT Int. Appl., 100 pp.

CODEN: PIXXD2

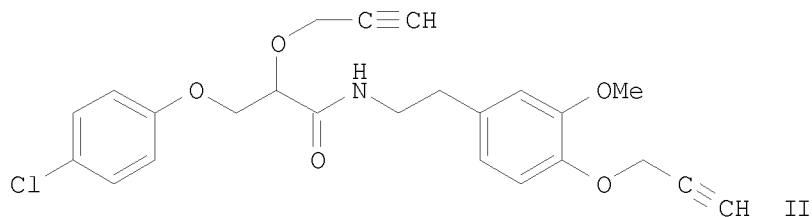
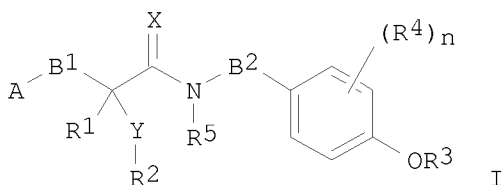
DOCUMENT TYPE: Patent

LANGUAGE: English

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003042167 A1		20030522	WO 2002-XB12845	20021115
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZM, ZW			
RW:	AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG, TR			
PRIORITY APPLN. INFO.:			GB 2001-27556	20011116

GI



AB Title compds. I [wherein A = (un)substituted (hetero)aryl; X = O or S; Y = O or S; R1 = H, (halo)alkyl, (halo)alkenyl, (halo)alkynyl, or (halo)cycloalkyl; R2 and R3 = independently H, (halo)alkyl, (halo)alkenyl, (halo)alkynyl, (halo)cycloalkyl(alkyl), (halo)alkoxyalkyl, (halo)alkoxyalkenyl, (halo)alkoxyalkynyl, or (un)substituted aryl(oxy)alkyl, arylalkenyl, or arylalkynyl; or R3 = (un)substituted

heteroarylalkyl, heteroarylalkenyl, or heteroarylalkynyl; R4 = halo, CN, NO2, NH2, CHO, CO2H, (halo)alkyl, (halo)alkenyl(oxy), (halo)alkynyl(oxy), (halo)alkoxy(alkyl), (halo)alkylthio, (halo)alkanoyl, (halo)(di)alkylamino, or (halo)alkoxycarbonyl; R5 = H, alkyl, alkenyl, or alkynyl; n = 0-4; B1 = (CR10R11)q or (CHR10R11)rZ(CR12R13)s; q = 2-4; r = 0-3; s = 1-3; r + s = 1-3; Z = O, S, SO, SO2, NR6, CO, OCO, CO2, NR6CO, or CONR6; R6 = H or alkyl; R10-R13 = independently H or alkyl; B2 = alkylene bridge; and optical isomers and mixts. thereof] were prepared These compds. possess useful plant protecting properties and may be employed advantageously in agricultural practice for controlling or preventing the infestation of plants by phytopathogenic microorganisms, especially fungi. For example, 3-chlorolactic acid was coupled with 4-chlorophenol in 3.3N NaOH to give 3-(4-chlorophenoxy)-2-hydroxypropionic acid. Amidation with 2-[3-methoxy-4-[(prop-2-ynyl)oxy]phenyl]ethylamine•HCl in the presence of N,N-diisopropylethylamine in DMF, followed by etherification with propargyl bromide in toluene provided the N-(phenethyl)- α -(propargyloxy)propionamide II. The latter showed residual protective action and residual curative action against fungal infestation by Plasmopara viticola on vines, Phytophthora on tomato plants, and Phytophthora on potato plants by 80-100% at 200 ppm. [This abstract record is one of 6 records for this document necessitated by the large number of index entries required to fully index the document and publication system constraints.]

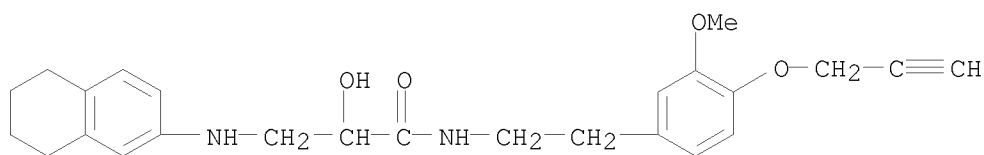
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RL: PRPH (Prophetic)

(Preparation of α -oxygenated or α -thiolated carboxylic acid phenethylamides for controlling fungal infestation in plants)

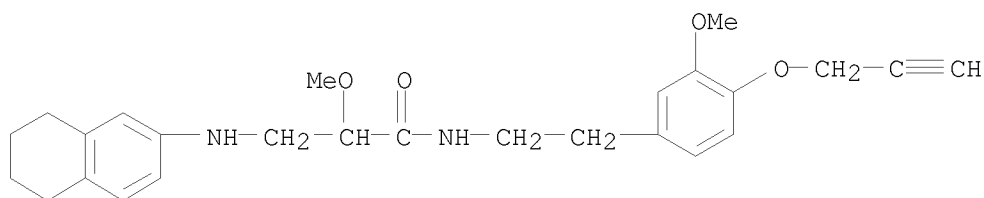
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RN 1067363-16-9 CAPLUS

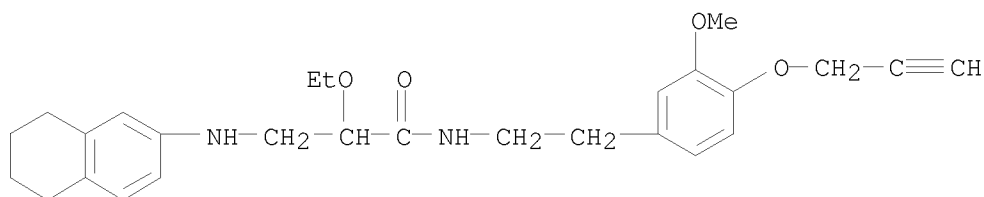
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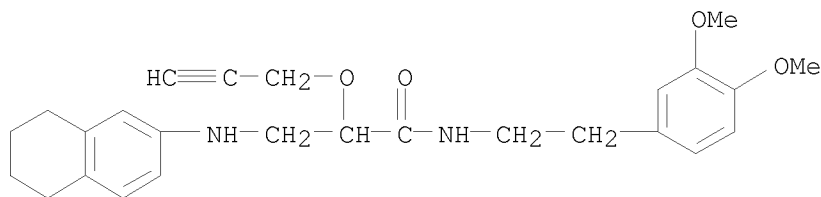
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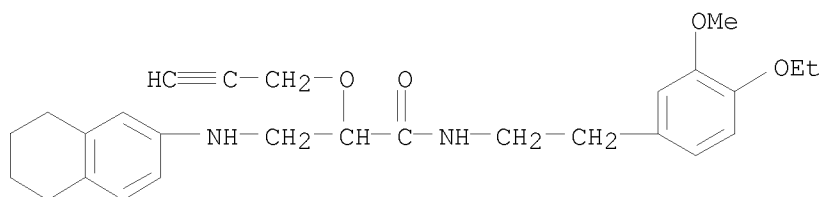
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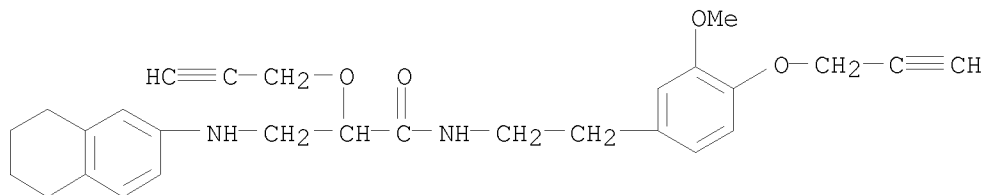
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RN 1067363-20-5 CAPLUS

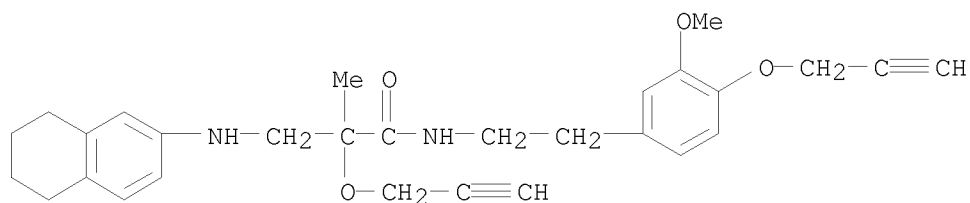
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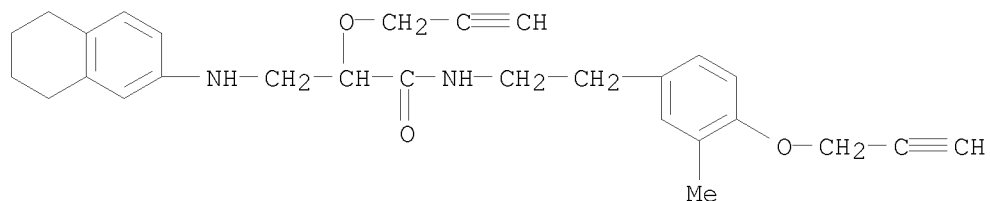
RN 1067363-21-6 CAPLUS

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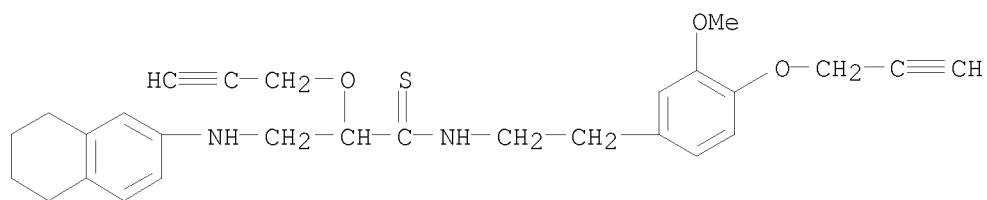
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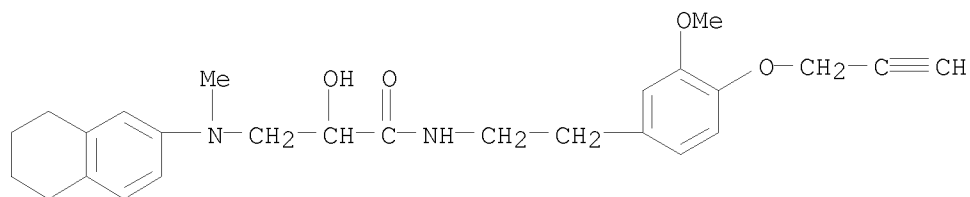
RN 1067363-23-8 CAPLUS

CN Propanethioamide, N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-2-(2-propyn-1-yloxy)-3-[(5,6,7,8-tetrahydro-2-naphthalenyl)amino]- (CA INDEX NAME)



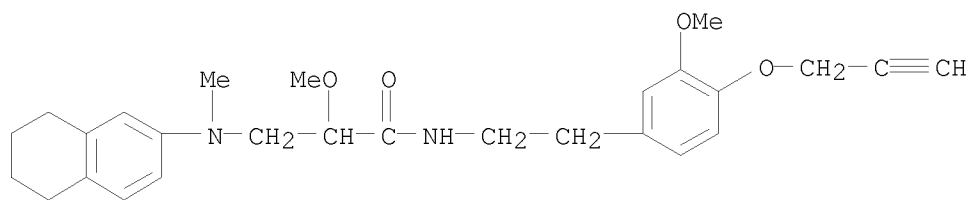
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CN INDEX NAME NOT YET ASSIGNED

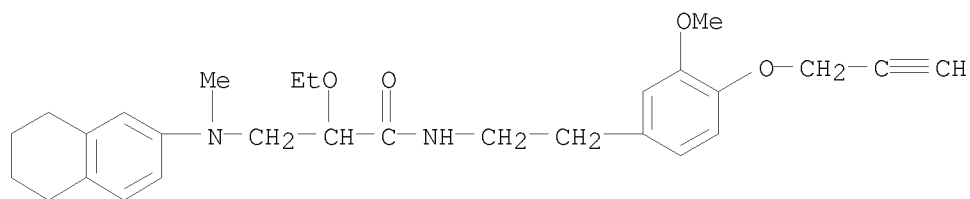


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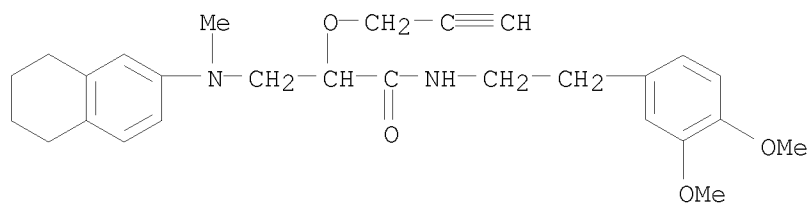
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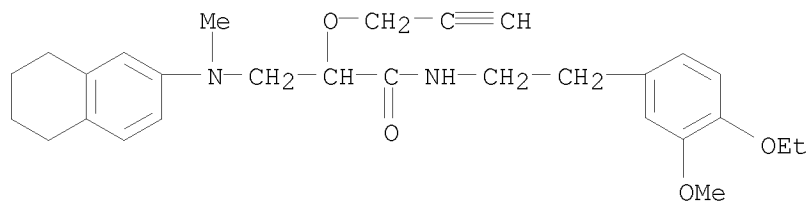
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CN INDEX NAME NOT YET ASSIGNED



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CN INDEX NAME NOT YET ASSIGNED

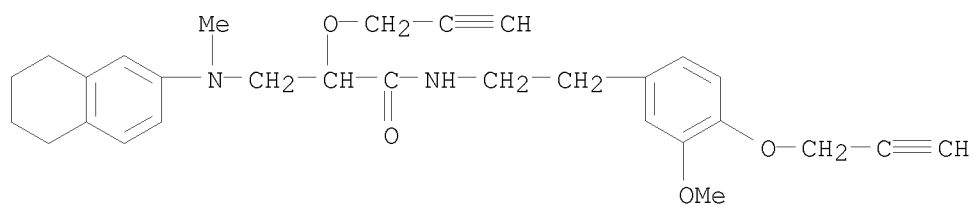


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CN INDEX NAME NOT YET ASSIGNED

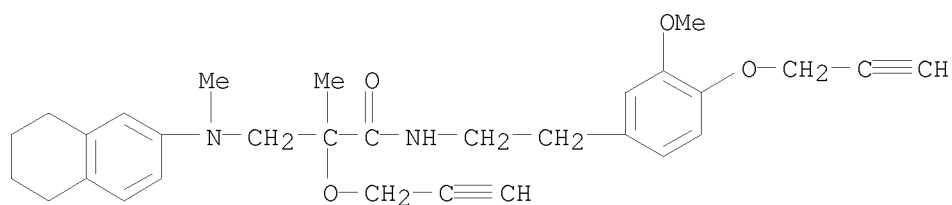


RN 1067407-86-6 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

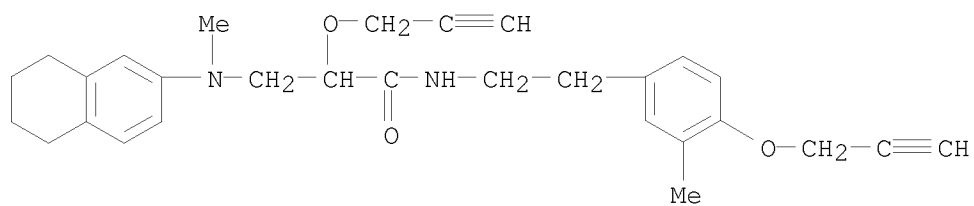
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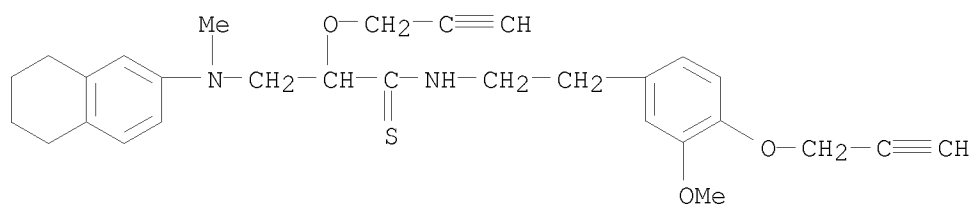
RN 1067407-87-7 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1067407-88-8 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1067407-89-9 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



REFERENCE COUNT: 7 THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L5 ANSWER 9 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2003:696854 CAPLUS

DOCUMENT NUMBER: 139:214722

TITLE: Preparation of (substituted)acyl dipeptidyl inhibitors of the ICE/ced-3 family of cysteine proteases

INVENTOR(S): Karanewsky, Donald S.; Kalish, Vincent J.; Robinson, Edard D.; Ullman, Brett R.

PATENT ASSIGNEE(S): Idun Pharmaceuticals, Inc., USA

SOURCE: PCT Int. Appl., 87 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003072528	A2	20030904	WO 2003-US3987	20030207 <--
WO 2003072528	A3	20040325		
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RW:	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
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AU 2003248360	A1	20030909	AU 2003-248360	20030207 <--
US 20030232788	A1	20031218	US 2003-360559	20030207 <--
EP 1480999	A2	20041201	EP 2003-743123	20030207 <--
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PRIORITY APPLN. INFO.:			US 2002-355390P	P 20020208
			WO 2003-US3987	W 20030207

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 139:214722

AB Compds. R1(CH₂)_nCHR₂CO-A-NHCH[(CH₂)_qCO₂R₃]CO-B [A is a natural or unnatural amino acid; B = H, D, alkyl, cycloalkyl, (un)substituted Ph or naphthyl, 2-benzoxazolyl, halomethyl, (CH₂)_mcycloalkyl, (CH₂)_m(1- or 2-naphthyl), substituted 2-oxazolyl, (un)substituted (CH₂)_mphenyl, CH₂OCO(aryl), or CH₂OCO(heteroaryl), etc.; R₁ = (un)substituted cycloalkyl, Ph, naphthyl, or heteroaryl; R₂ = H, alkyl, cycloalkyl, (un)substituted Ph, (CH₂)_mNH₂, (un)substituted (CH₂)_mphenyl, (CH₂)_mcycloalkyl, (CH₂)_mheteroaryl, etc.; R₃ = H, alkyl, cycloalkyl, (cycloalkyl)alkyl, (un)substituted phenylalkyl; m = 1-4, n = 0-2; q = 1-2] or their pharmaceutically-acceptable salts were prepared as inhibitors of the ICE/ced-3 family of cysteine proteases (ICE = interleukin-1 β converting enzyme). Thus, coupling of (1-naphthyl)acetic acid with (3S)-3-(leucinylamino)-4-oxobutanoic acid tert-Bu ester semicarbazone (preparation given) followed by deprotection of the resulting intermediate with TFA, and treatment with a 3:1:1 solution of MeOH/AcOH/37% HCHO afforded (3S)-3-[[N-[(1-naphthyl)acetyl]leuciny]amino]-4-oxobutanoic acid. The invention is also directed to pharmaceutical compns. containing these compds., as well as the use of such compns. in the treatment of patients suffering

10/513699

inflammatory, autoimmune and neurodegenerative diseases, for the prevention of ischemic injury, and for the preservation of organs that are to undergo a transplantation procedure.

IT 1080814-59-0

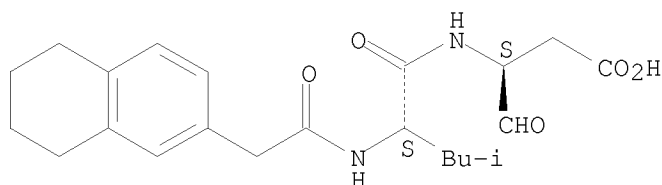
RL: PRPH (Prophetic)

(Preparation of (substituted)acyl dipeptidyl inhibitors of the ICE/ced-3 family of cysteine proteases)

RN 1080814-59-0 CAPLUS

CN INDEX NAME NOT YET ASSIGNED

Absolute stereochemistry.



OS.CITING REF COUNT:	1	THERE ARE 1 CAPLUS RECORDS THAT CITE THIS RECORD (1 CITINGS)
REFERENCE COUNT:	3	THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L5 ANSWER 10 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2003:472358 CAPLUS

DOCUMENT NUMBER: 139:53025

TITLE: Preparation of vanilloid receptor ligands and their use in treatments

INVENTOR(S): Bo, Yunxin Y.; Chakrabarti, Partha P.; Chen, Ning; Doherty, Elizabeth M.; Fotsch, Christopher H.; Han, Nianhe; Kelly, Michael G.; Liu, Qingyian; Norman, Mark Henry; Wang, Xianghong; Zhu, Jiawang; Ognyanov, Vassil; Bo, Yunxin Y.; Chakrabarti, Partha P.; Chen, Ning; Doherty, Elizabeth M.; Fotsch, Christopher H.; Han, Nianhe; Kelly, Michael; Liu, Qingyian; et al.

PATENT ASSIGNEE(S): Amgen Inc., USA; et al.

SOURCE: PCT Int. Appl., 611 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003049702	A2	20030619	WO 2002-US39589	20021210 <--
WO 2003049702	A3	20040212		
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AU 2002364549	B2	20071122		
US 20030195201	A1	20031016	US 2002-316295	20021210 <--
US 7582657	B2	20090901		
EP 1463714	A2	20041006	EP 2002-799927	20021210 <--
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EP 1764358	A2	20070321	EP 2006-10087	20021210
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EP 1780196	A2	20070502	EP 2006-10095	20021210
EP 1780196	A3	20070509		
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WO 2003099284	A1	20031204	WO 2003-US16655	20030520 <--
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AU 2003247425	A1	20031212	AU 2003-247425	20030520 <--
AU 2003247425	B2	20070308		
US 20040038969	A1	20040226	US 2003-445170	20030520 <--
US 7053088	B2	20060530		
EP 1542692	A1	20050622	EP 2003-755509	20030520
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JP 2005531574	T	20051020	JP 2004-506808	20030520
EP 1688408	A2	20060809	EP 2006-8551	20030808
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MX 2004005427	A	20050419	MX 2004-5427	20040604
MX 2004011472	A	20050214	MX 2004-11472	20041118
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US 7579347	B2	20090825		
US 20050272931	A1	20051208	US 2005-99978	20050405
US 20060030618	A1	20060209	US 2005-100272	20050405
US 20050267163	A1	20051201	US 2005-195302	20050801
US 7524874	B2	20090428		
US 20050272777	A1	20051208	US 2005-195159	20050801
US 7332511	B2	20080219		
US 20050277631	A1	20051215	US 2005-195134	20050801
US 7148221	B2	20061212		
US 20050277646	A1	20051215	US 2005-195303	20050801
US 7396831	B2	20080708		
AU 2007200149	A1	20070201	AU 2007-200149	20070115
AU 2008202517	A1	20080626	AU 2008-202517	20080605
US 20090264424	A1	20091022	US 2009-492376	20090626
PRIORITY APPLN. INFO.:				
			US 2001-339161P	P 20011210
			US 2001-344737P	P 20011221
			US 2002-383331P	P 20020522
			US 2002-402422P	P 20020808
			AU 2002-364549	A3 20021210
			EP 2002-799927	A3 20021210
			US 2002-316295	A3 20021210
			WO 2002-US39589	W 20021210
			US 2003-445170	A3 20030520
			WO 2003-US16655	W 20030520
			AU 2003-264047	A3 20030808
			EP 2003-785220	A3 20030808
			US 2003-638009	A3 20030808
			US 2005-100077	A3 20050405

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 139:53025

AB Claimed are compds. having the general structure R1CR2:CR3C(:X)YR4 or
 R1R2CHCR3R3C(:X)YR4 (I; variables defined below; e.g.

(2E)-3-[4-(tert-butyl)phenyl]-N-phenylprop-2-enamide and (2,3-dihydrobenzo[1,4]dioxin-6-yl)[4-(4-dimethylaminophenyl)pyridin-2-yl]amine) and compns. containing them, for the treatment of acute, inflammatory and neuropathic pain, dental pain, general headache, migraine, cluster headache, mixed-vascular and nonvascular syndromes, tension headache, , general inflammation arthritis, rheumatic diseases, osteoarthritis, inflammatory bowel disorders, inflammatory eye disorders, inflammatory or unstable bladder disorders, psoriasis, skin complaints with inflammatory components, chronic inflammatory conditions, inflammatory pain and associated hyperalgesia and allodynia, neuropathy pain and associated hyperalgesia and allodynia, diabetic neuropathy pain, causalgia, sympathetically maintained pain, deafferentiation syndromes, asthma, epithelial tissue damage or dysfunction, herpes simplex, disturbances of visceral motility at respiratory, genitourinary, gastrointestinal or vascular regions, wounds, burns, allergic skin reactions, pruritis, vitiligo, general gastrointestinal disorders, gastric ulceration, duodenal ulcers, diarrhea, gastric lesions induced by necrotising agents, hair growth, vasomotor or allergic rhinitis, bronchial disorders or bladder disorders. I are thought to be vanilloid receptor ligands, but no test data are provided. Although the methods of preparation are not claimed, .apprx.130 example prepns. and characterization data for .apprx.400 I are included. For I: R1 is Ph, naphthyl or (un)saturated 5- or 6-membered ring heterocycle; R2 is H, hydroxy, halo, C1-6alkyl, or (un)saturated 5- or 6-membered ring heterocycle; or R1 and R2 together are o-benzenediyl-L1-o-benzenediyl. R3 is H or C1-4alkyl; or R1 and R3 together are o-benzenediyl-L2- or -Z-L2- (Z = pyridine-2,3-diyl). R4 is Ph, (un)saturated 5- or 6-membered ring heterocycle, 10-membered bicyclic ring comprising fused 6-membered rings, containing 0-4 N atoms with the remainder being C atoms, with at least one of the 6-membered rings being aromatic; X is O, S or NRa; or X and R2 together are :N-CH:CH-, :C-O-, :C-S-, or :C-NRa-; Y is NH or O; addnl. details including provisos are given in the claims.

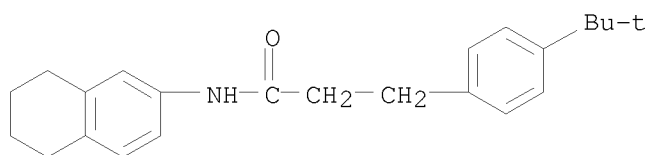
IT 1064718-76-8

RL: PRPH (Prophetic)

(Preparation of vanilloid receptor ligands and their use in treatments)

RN 1064718-76-8 CAPLUS

CN Benzenepropanamide, 4-(1,1-dimethylethyl)-N-(5,6,7,8-tetrahydro-2-naphthalenyl)- (CA INDEX NAME)



OS.CITING REF COUNT: 29 THERE ARE 29 CAPLUS RECORDS THAT CITE THIS RECORD (66 CITINGS)

10/513699

L5 ANSWER 11 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2003:396732 CAPLUS

DOCUMENT NUMBER: 138:385175

TITLE: Preparation of
N-[[(propargyloxy)phenyl]alkyl]arylacetamides for
controlling fungal infestations in plants

INVENTOR(S): Zeller, Martin; Lamberth, Clemens

PATENT ASSIGNEE(S): Syngenta Participations A.-G., Switz.

SOURCE: PCT Int. Appl., 61 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

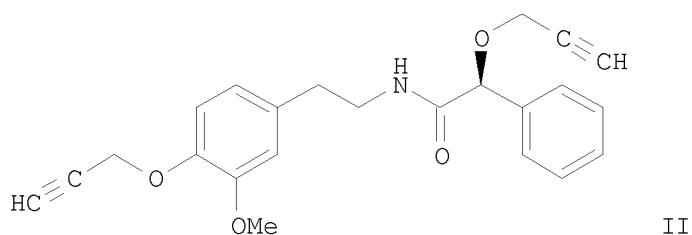
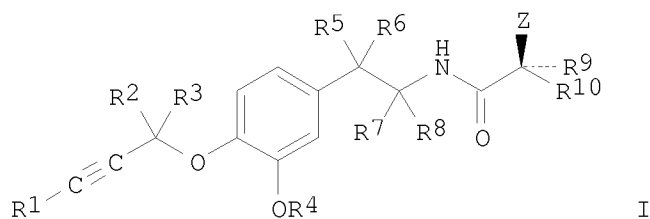
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2003041728	A1	20030522	WO 2002-EP12848	20021115 <--
WO 2003041728	A9	20040422		
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CA 2464807	A1	20030522	CA 2002-2464807	20021115 <--
AU 2002363600	A1	20030526	AU 2002-363600	20021115 <--
EP 1444197	A1	20040811	EP 2002-798312	20021115 <--
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BR 2002014139	A	20041019	BR 2002-14139	20021115 <--
AT 292108	T	20050415	AT 2002-798312	20021115
JP 2005519030	T	20050630	JP 2003-543615	20021115
ES 2239272	T3	20050916	ES 2002-798312	20021115
IN 2004DN01098	A	20050401	IN 2004-DN1098	20040423
US 20050026785	A1	20050203	US 2004-495153	20040510
MX 2004004550	A	20040813	MX 2004-4550	20040513 <--
CN 101421230	A	20090429	CN 2002-822692	20040514
KR 897729	B1	20090515	KR 2004-707452	20040514
PRIORITY APPLN. INFO.:			GB 2001-27554	A 20011116
			WO 2002-EP12848	W 20021115

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 138:385175

GI



AB Title Ph propargyl ether derivs. I [wherein R1 = H, (cyclo)alkyl, or (un)substituted aryl; R2 and R3 = independently H or alkyl; R4 = aryl, alkenyl, or alkynyl; R5-R8 = independently H or alkyl; R9 = H or (un)substituted alkyl, alkenyl, or alkynyl; R10 = (un)substituted (hetero)aryl; Z = (un)substituted aryloxy, alkoxy, alkenyloxy, or alkynyloxy; and optical isomers and mixts. thereof] were prepared These compds. possess useful plant protecting properties and may be employed advantageously in agricultural practice for controlling or preventing the infestation of plants by phytopathogenic microorganisms, especially fungi. For example, reaction of 2-[3-methoxy-4-[(prop-2-ynyl)oxy]phenyl]ethylamine•HCl and L-(+)-mandelic acid in the presence of N,N-diisopropylethylamine in DMF gave the amide. Etherification with propargyl bromide in toluene provided II. The latter showed residual protective action and residual curative action against fungal infestation by *Plasmopara viticola* on vines, *Phytophthora* on tomato plants, and *Phytophthora* on potato plants by 80-100% at 200 ppm.

IT	1055179-98-0	1055181-48-0	1055182-12-1
	1055182-57-4	1055182-58-5	1055182-59-6
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1055226-57-7	1055228-55-1	1055228-56-2
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1055230-42-6	1055230-43-7	1055230-44-8
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1055232-71-7	1055236-78-6	1055238-63-5
1055238-64-6	1055238-65-7	1055238-66-8
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1055243-87-2	1055243-88-3	1055243-89-4
1055244-11-5	1055244-12-6	1055244-13-7
1055244-14-8	1055244-15-9	1055247-12-5
1055247-15-8	1055247-16-9	1055249-28-9
1055249-29-0	1055249-30-3	1055249-33-6
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1055250-45-7	1055253-12-7	1055254-60-8
1055254-61-9	1055254-62-0	1055254-63-1
1055254-64-2	1055255-69-0	1055255-70-3
1055255-71-4	1055255-72-5	1055255-73-6
1055256-73-9	1055258-31-5	1055258-32-6
1055258-33-7	1055261-20-5	1055261-21-6
1055261-22-7	1055261-23-8	1055261-24-9
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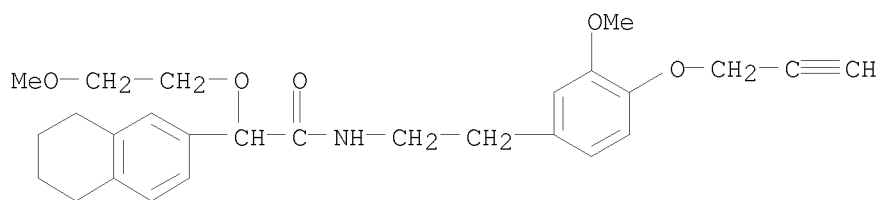
RL: PRPH (Prophetic)

(Preparation of N-[[(propargyloxy)phenyl]alkyl]arylacetamides for
controlling fungal infestations in plants)

RN 1055179-98-0 CAPLUS

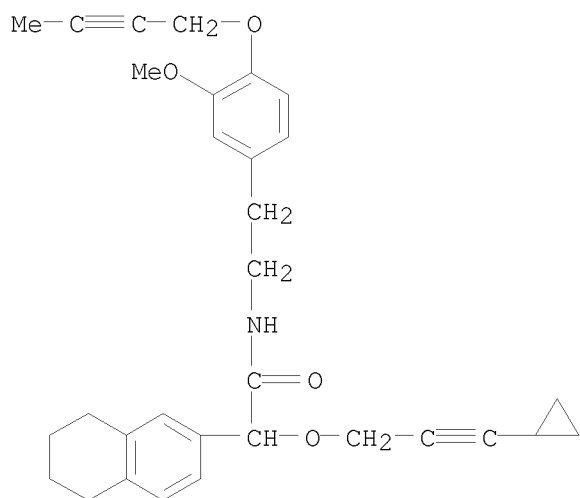
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -(2-methoxyethoxy)-N-[2-
[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

10/513699



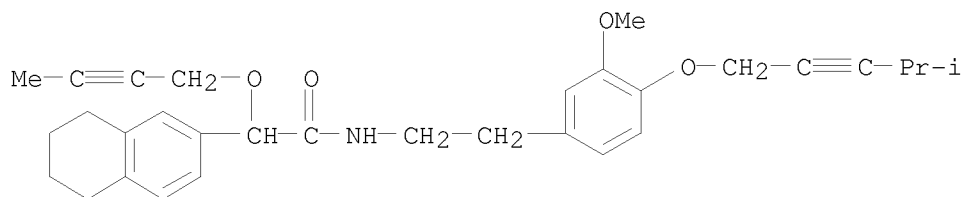
RN 1055181-48-0 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-
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NAME)



RN 1055182-12-1 CAPLUS

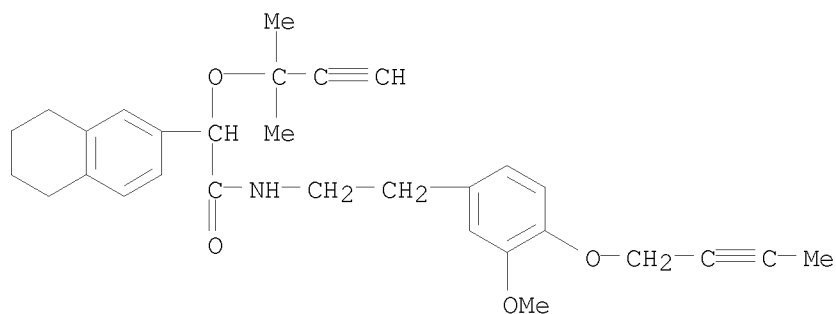
CN INDEX NAME NOT YET ASSIGNED



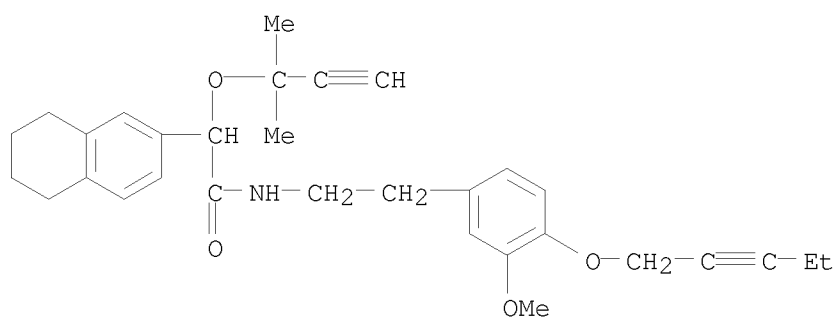
RN 1055182-57-4 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-
α-[(1,1-dimethyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro- (CA INDEX
NAME)

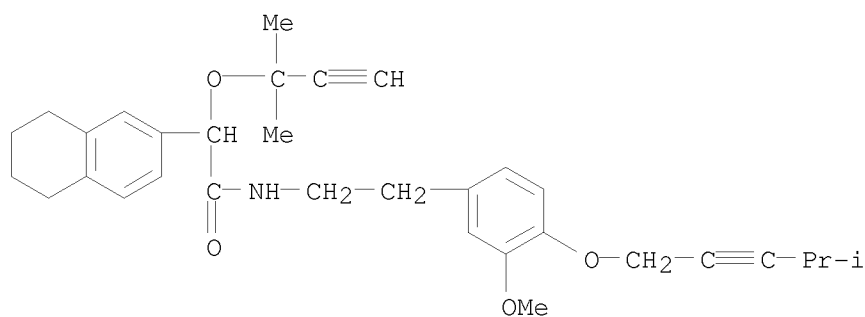
10/513699



RN 1055182-58-5 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

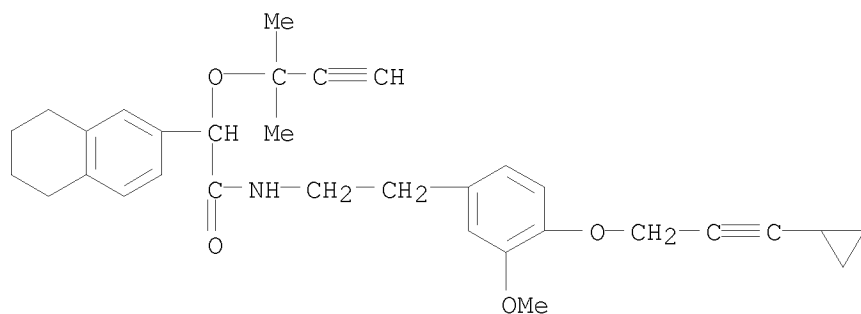


RN 1055182-59-6 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



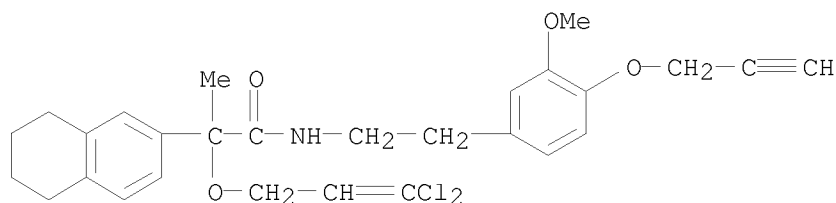
RN 1055182-60-9 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-alpha-[(1,1-dimethyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro- (CA INDEX NAME)

10/513699



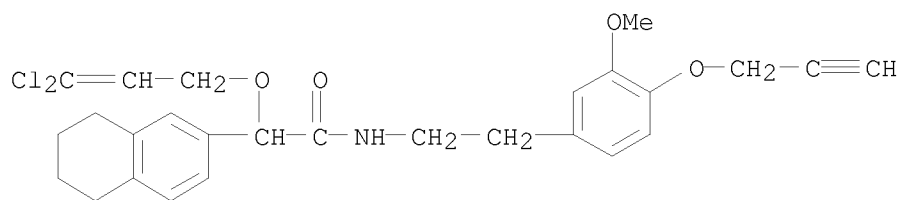
RN 1055183-36-2 CAPLUS

CN 2-Naphthaleneacetamide, α -[(3,3-dichloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



RN 1055183-37-3 CAPLUS

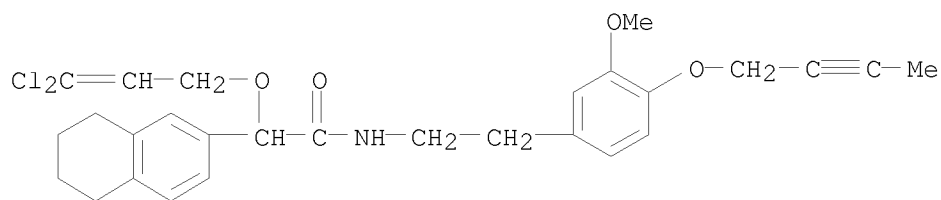
CN 2-Naphthaleneacetamide, α -[(3,3-dichloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



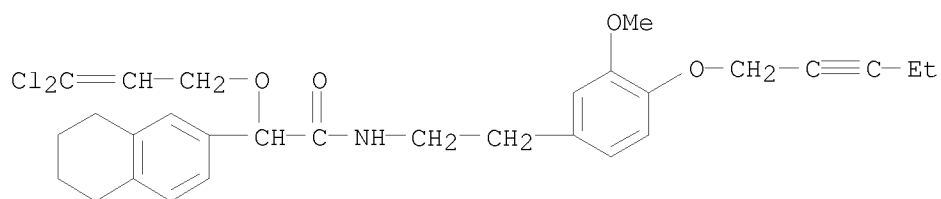
RN 1055183-38-4 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -[(3,3-dichloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro- (CA INDEX NAME)

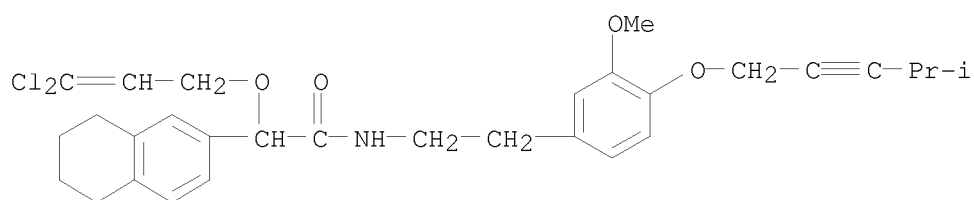
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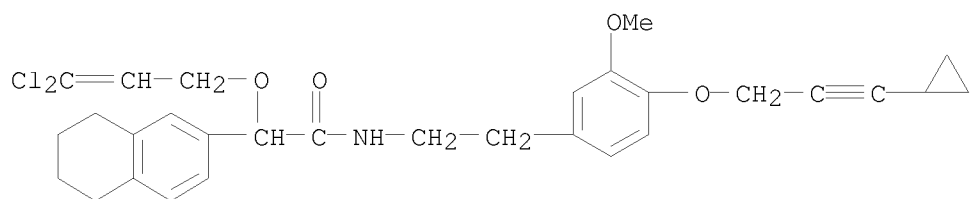
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RN 1055183-40-8 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

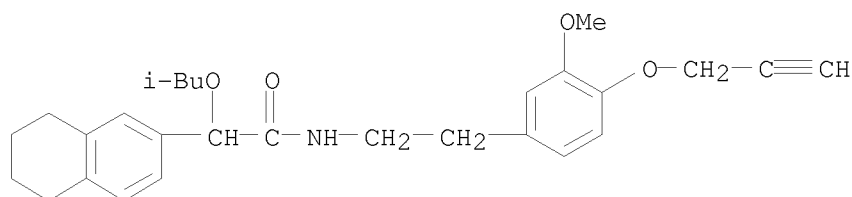


RN 1055183-41-9 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

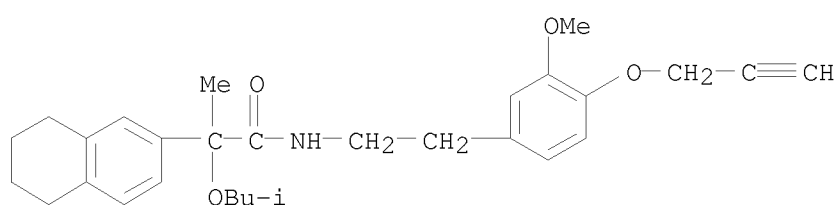


RN 1055185-17-5 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(2-methylpropoxy)- (CA INDEX NAME)

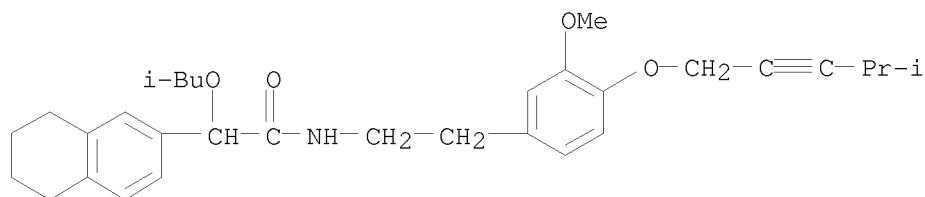
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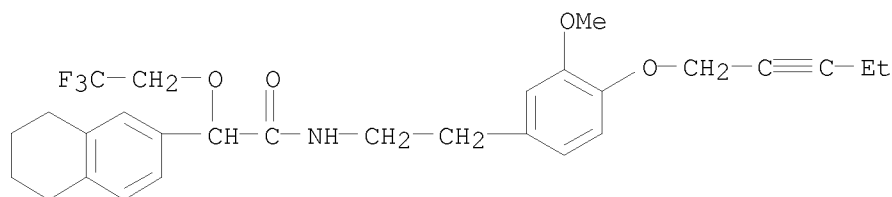
RN 1055185-18-6 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-(2-methylpropoxy)- (CA INDEX NAME)



RN 1055185-19-7 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

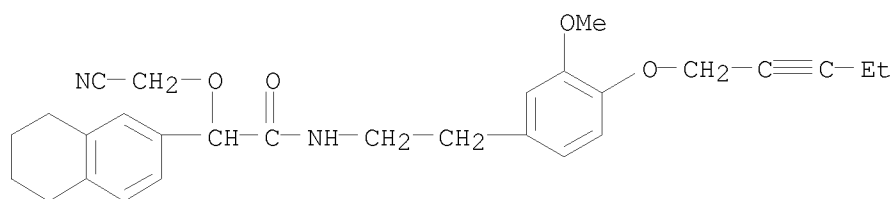


RN 1055186-14-5 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]-α-(2,2,2-trifluoroethoxy)- (CA INDEX NAME)

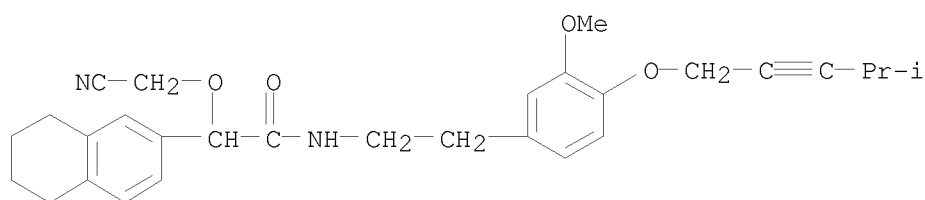


RN 1055186-58-7 CAPLUS
CN 2-Naphthaleneacetamide, α-(cyanomethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

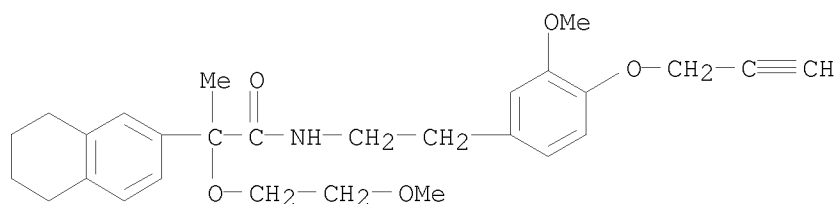
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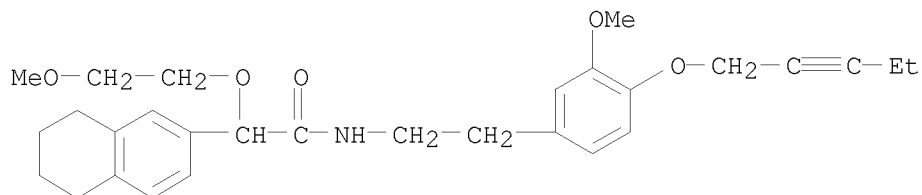
RN 1055186-59-8 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055189-32-6 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -(2-methoxyethoxy)-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)

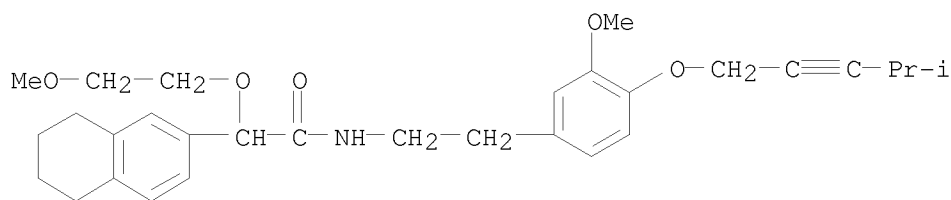


RN 1055189-33-7 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -(2-methoxyethoxy)-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



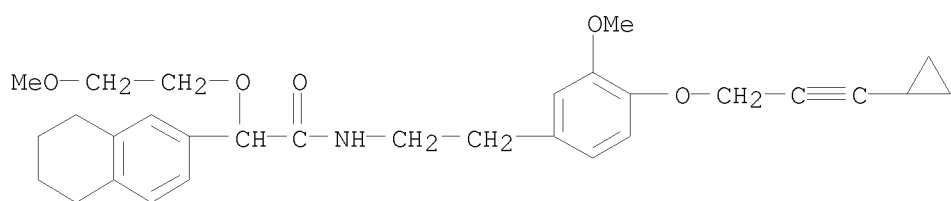
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CN INDEX NAME NOT YET ASSIGNED

10/513699



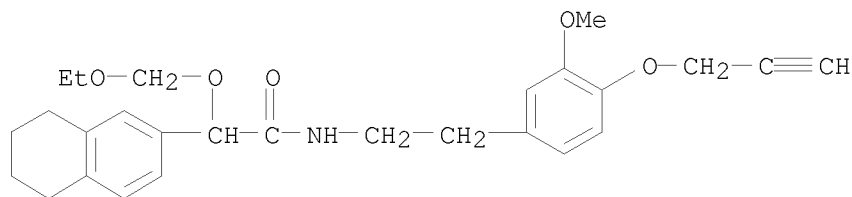
RN 1055189-35-9 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-methoxyethoxy)- (CA INDEX NAME)



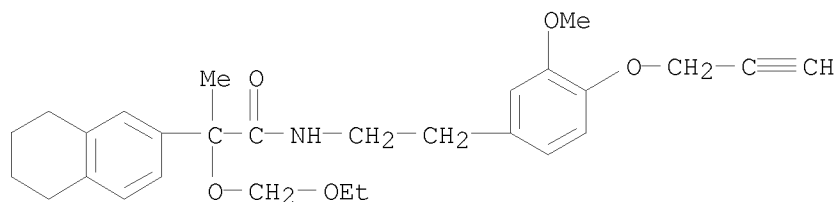
RN 1055199-07-9 CAPLUS

CN 2-Naphthaleneacetamide, α -(ethoxymethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055199-08-0 CAPLUS

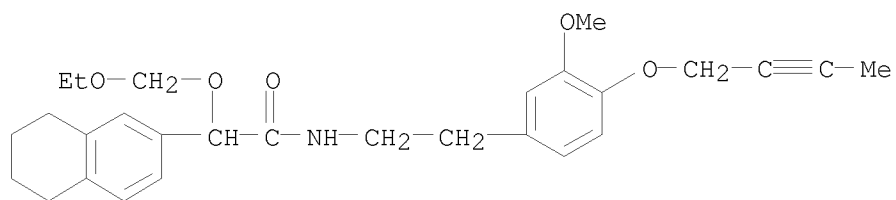
CN 2-Naphthaleneacetamide, α -(ethoxymethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



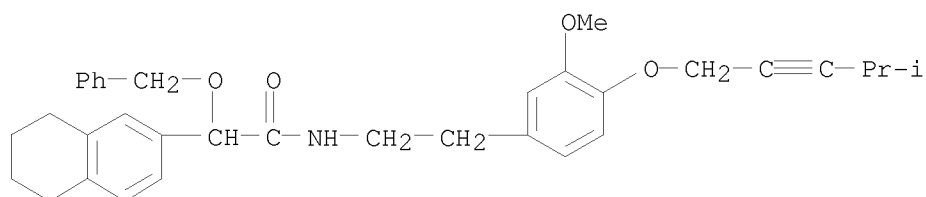
RN 1055199-09-1 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -(ethoxymethoxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)

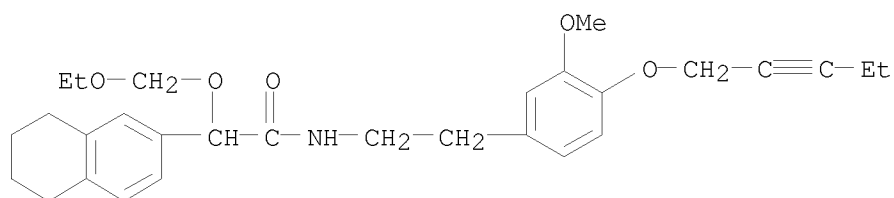
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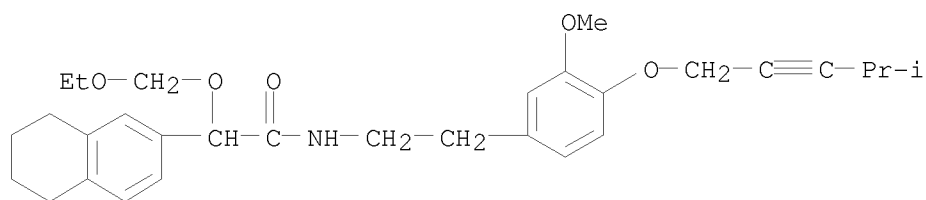
RN 1055203-51-4 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055203-63-8 CAPLUS
CN 2-Naphthaleneacetamide, α -(ethoxymethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

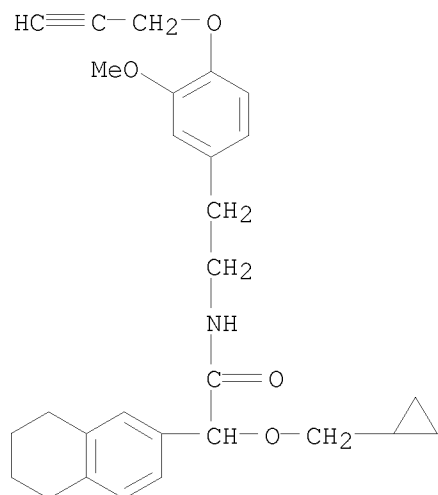


RN 1055203-64-9 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



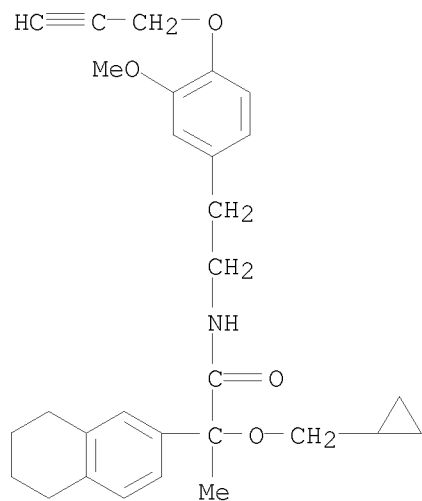
RN 1055206-24-0 CAPLUS
CN 2-Naphthaleneacetamide, α -(cyclopropylmethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

10/513699



RN 1055206-25-1 CAPLUS

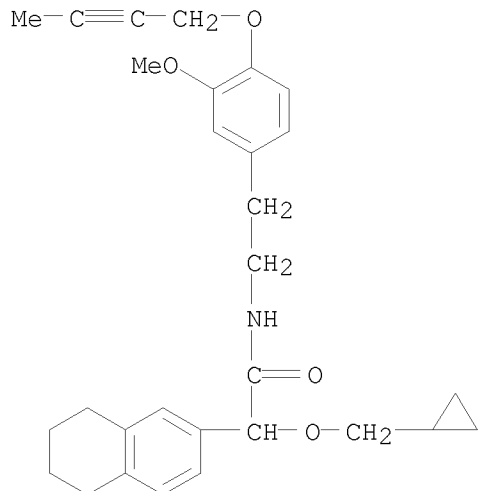
CN 2-Naphthaleneacetamide, α -(cyclopropylmethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



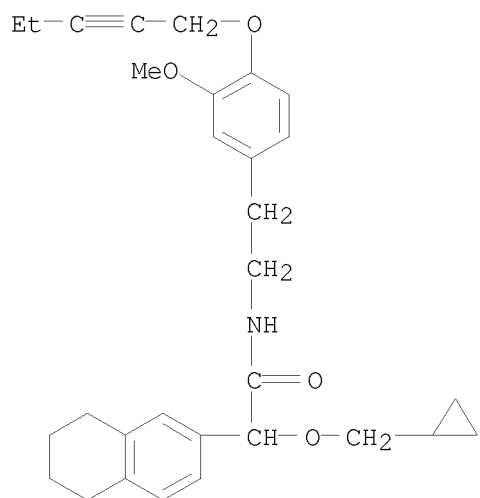
RN 1055206-26-2 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -(cyclopropylmethoxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)

10/513699

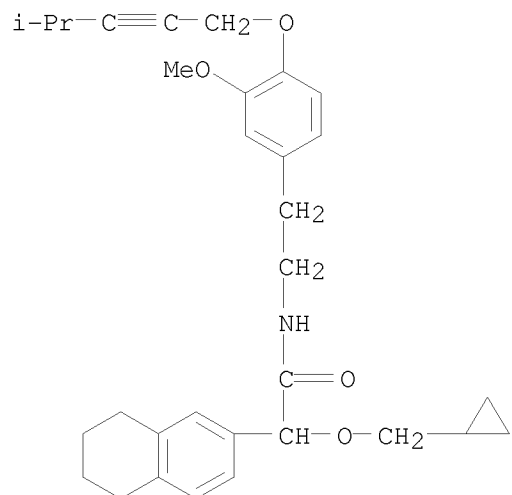


RN 1055206-27-3 CAPLUS
CN 2-Naphthaleneacetamide, α -(cyclopropylmethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



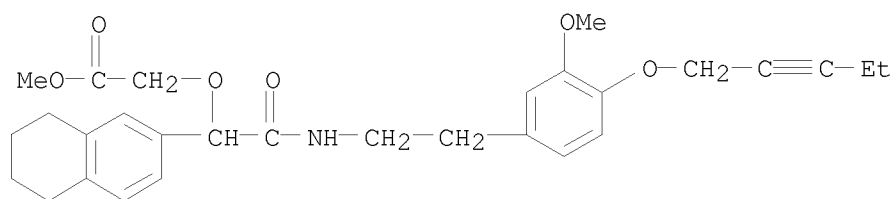
RN 1055206-28-4 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

10/513699



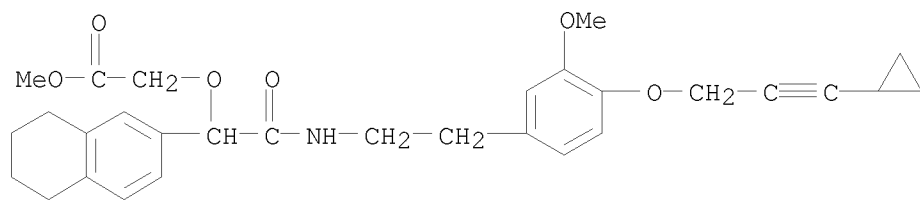
RN 1055207-50-5 CAPLUS

CN Acetic acid, 2-[2-[[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]amino]-2-oxo-1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethoxy]-, methyl ester (CA INDEX NAME)



RN 1055207-51-6 CAPLUS

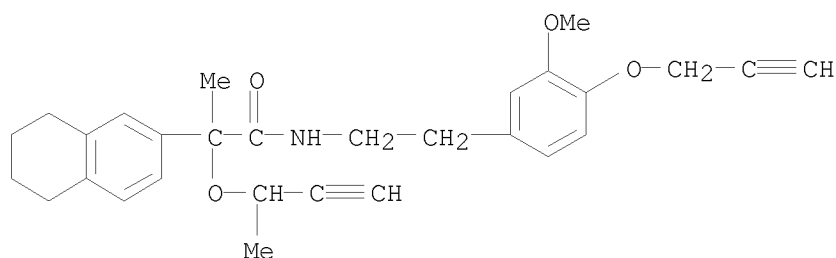
CN Acetic acid, 2-[2-[[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]amino]-2-oxo-1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethoxy]-, methyl ester (CA INDEX NAME)



RN 1055208-05-3 CAPLUS

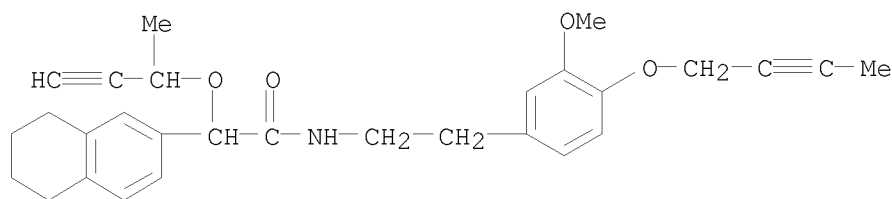
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-[(1-methyl-2-propyn-1-yl)oxy]- (CA INDEX NAME)

10/513699



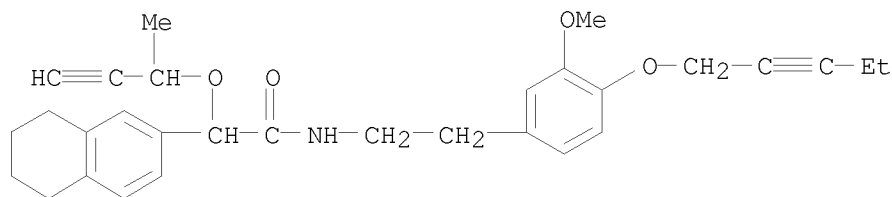
RN 1055208-06-4 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-[(1-methyl-2-propyn-1-yl)oxy]- (CA INDEX NAME)



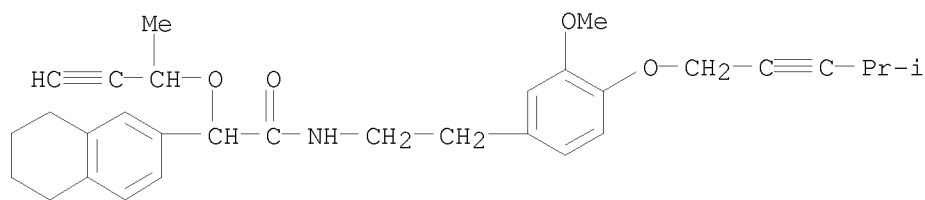
RN 1055208-07-5 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



RN 1055208-08-6 CAPLUS

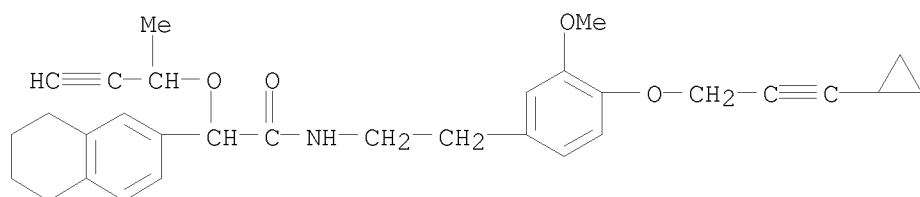
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RN 1055208-09-7 CAPLUS

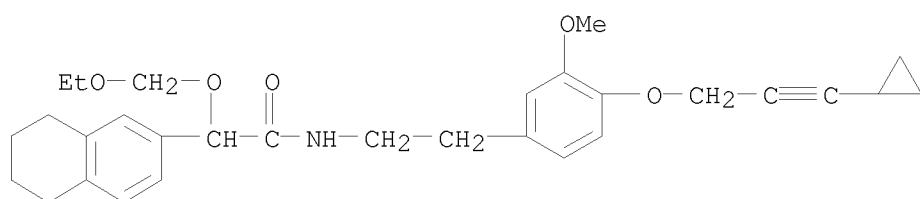
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-[(1-methyl-2-propyn-1-yl)oxy]- (CA INDEX NAME)

10/513699



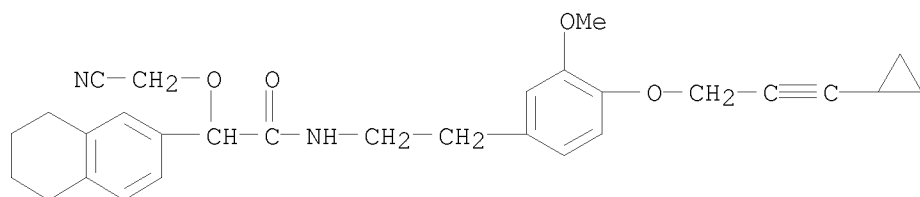
RN 1055208-72-4 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-α-(ethoxymethoxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)



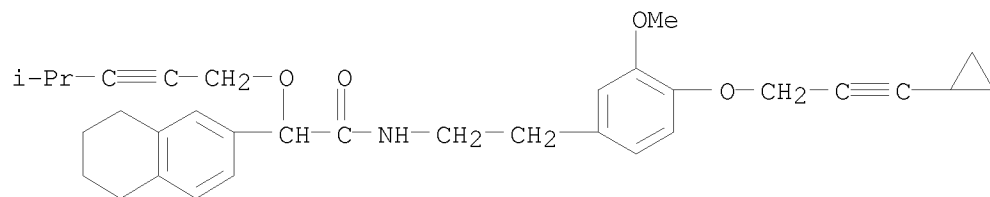
RN 1055209-45-4 CAPLUS

CN 2-Naphthaleneacetamide, α-(cyanomethoxy)-N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 1055210-27-9 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



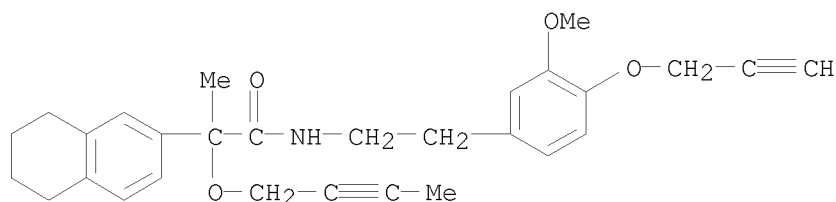
RN 1055214-37-3 CAPLUS

CN 2-Naphthaleneacetamide, α-(2-butyn-1-yloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl- (CA INDEX NAME)

<12/04/2007>

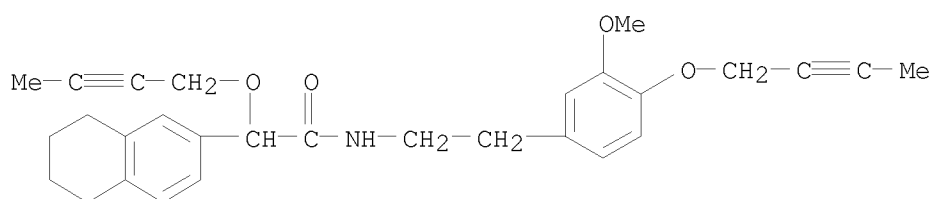
Erich Leese

10/513699



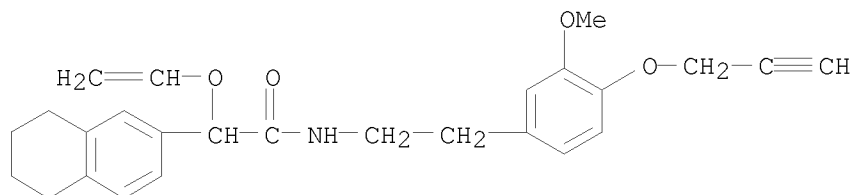
RN 1055214-38-4 CAPLUS

CN 2-Naphthaleneacetamide, α -(2-butyn-1-yloxy)-N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)



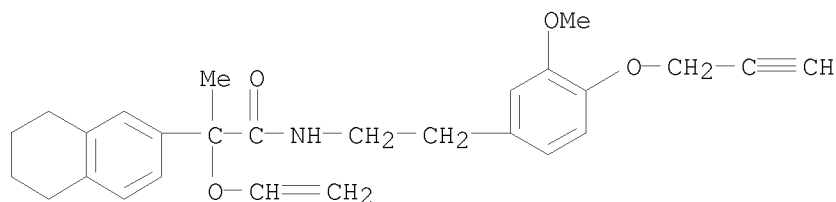
RN 1055215-71-8 CAPLUS

CN 2-Naphthaleneacetamide, α -(ethenyloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055215-72-9 CAPLUS

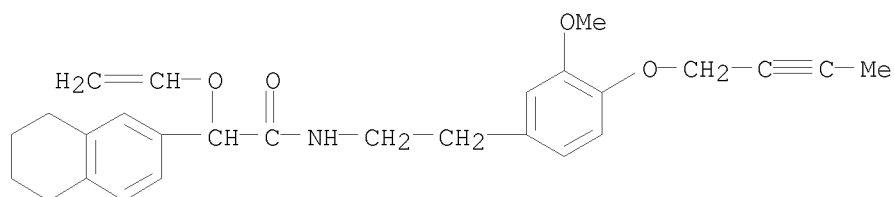
CN 2-Naphthaleneacetamide, α -(ethenyloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



RN 1055215-73-0 CAPLUS

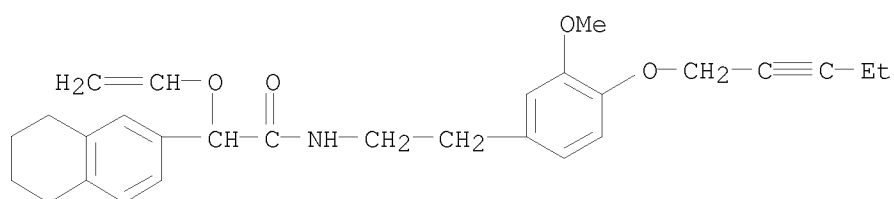
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -(ethenyloxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)

10/513699



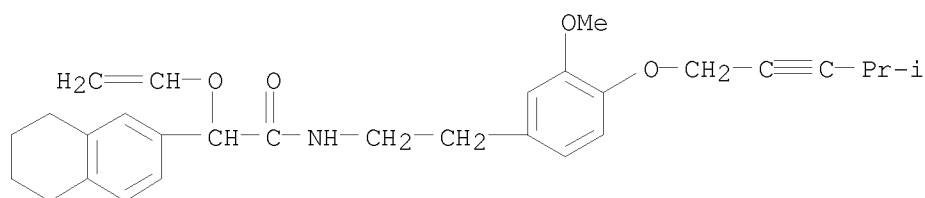
RN 1055215-74-1 CAPLUS

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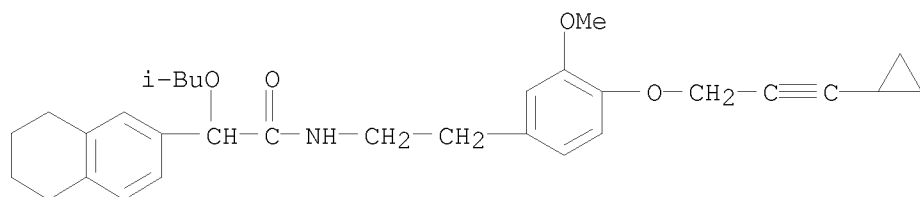
RN 1055215-75-2 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



RN 1055216-06-2 CAPLUS

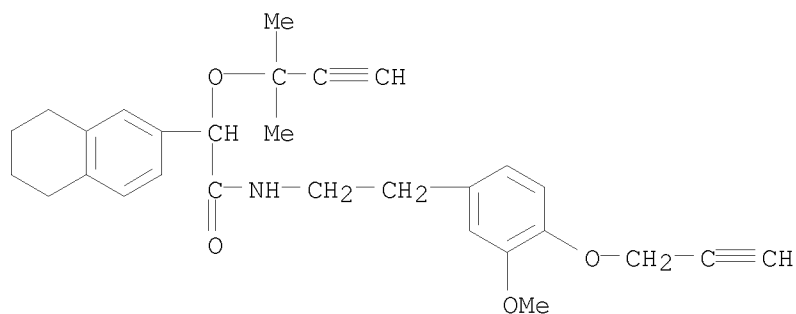
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-methylpropoxy)- (CA INDEX NAME)



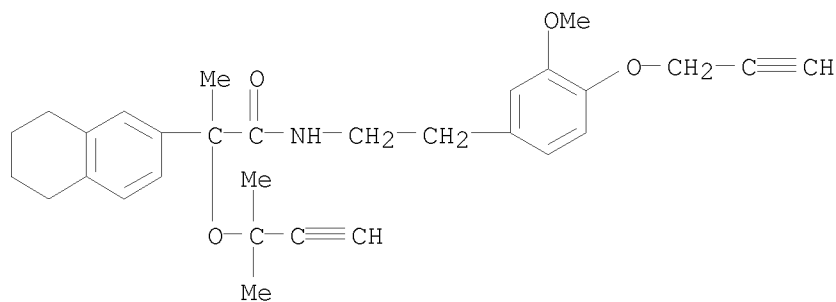
RN 1055216-39-1 CAPLUS

CN 2-Naphthaleneacetamide, α -[(1,1-dimethyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

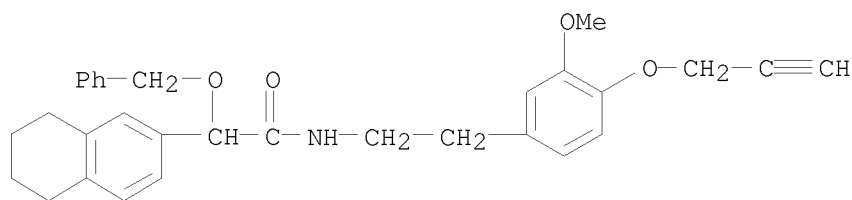
10/513699



RN 1055216-40-4 CAPLUS
CN 2-Naphthaleneacetamide, α -[(1,1-dimethyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)

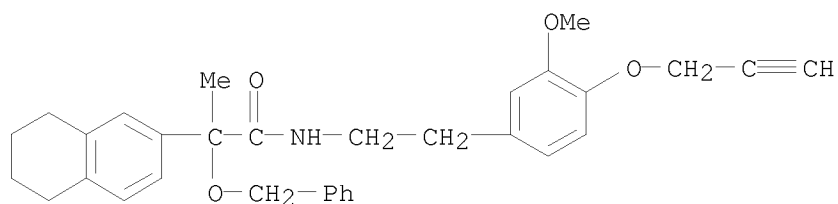


RN 1055216-75-5 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(phenylmethoxy)- (CA INDEX NAME)



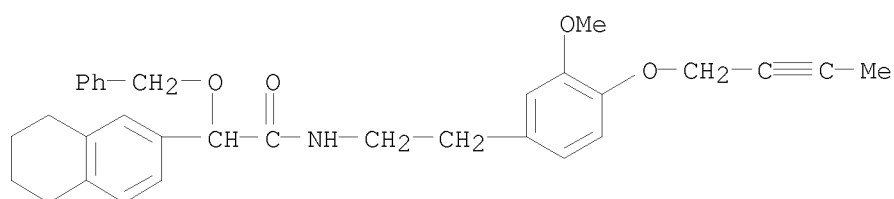
RN 1055216-76-6 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- α -(phenylmethoxy)- (CA INDEX NAME)

10/513699



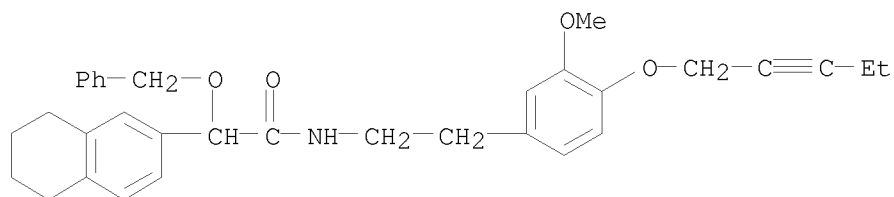
RN 1055216-77-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(phenylmethoxy)- (CA INDEX NAME)



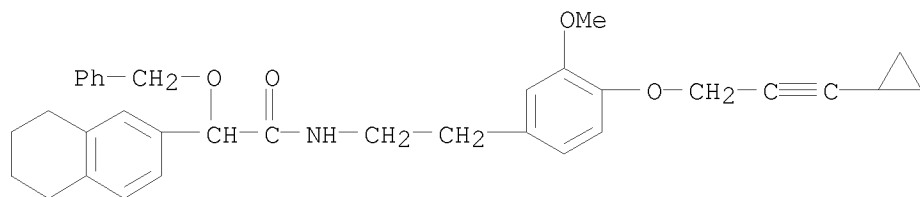
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CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- α -(phenylmethoxy)- (CA INDEX NAME)



RN 1055216-79-9 CAPLUS

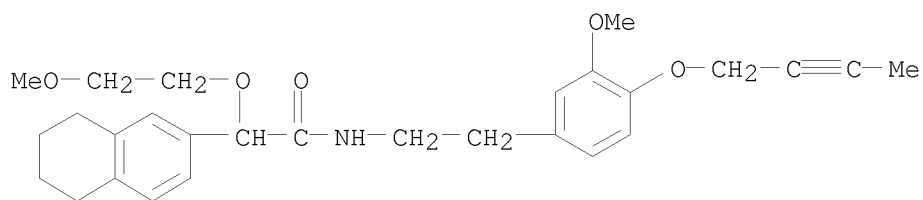
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(phenylmethoxy)- (CA INDEX NAME)



RN 1055216-94-8 CAPLUS

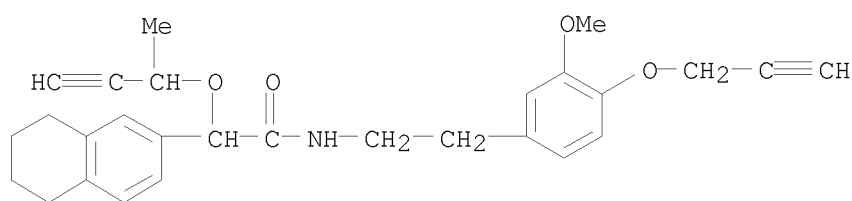
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-methoxyethoxy)- (CA INDEX NAME)

10/513699



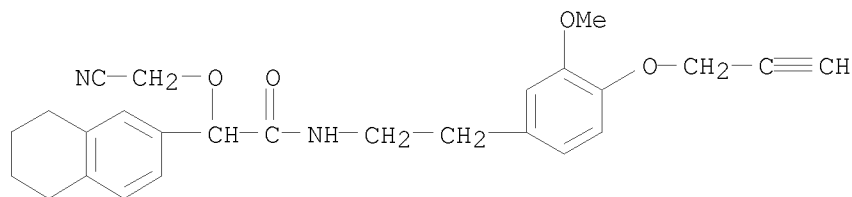
RN 1055217-05-4 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-[(1-methyl-2-propyn-1-yl)oxy]- (CA INDEX NAME)



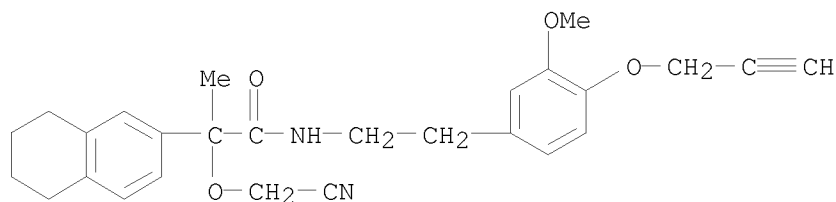
RN 1055217-88-3 CAPLUS

CN 2-Naphthaleneacetamide, α-(cyanomethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055217-89-4 CAPLUS

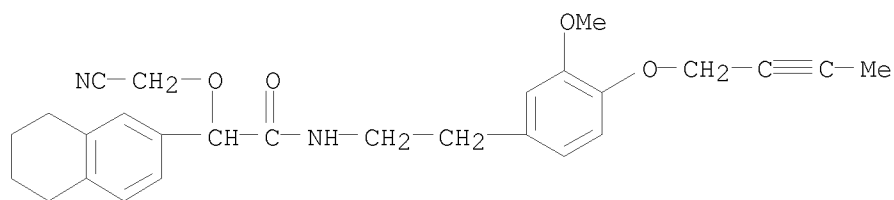
CN 2-Naphthaleneacetamide, α-(cyanomethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl- (CA INDEX NAME)



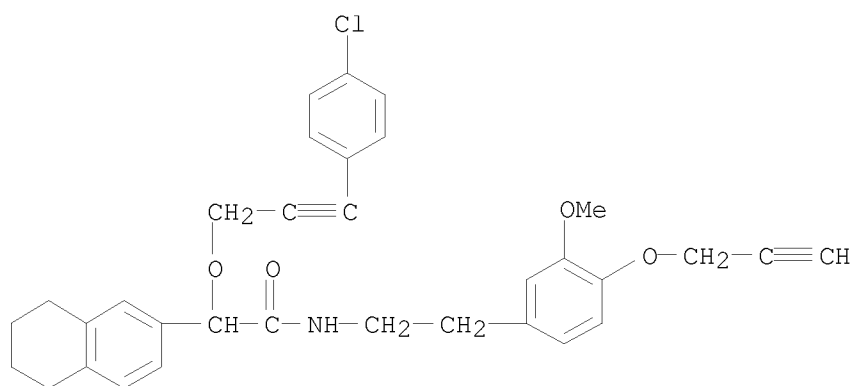
RN 1055217-90-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-α-(cyanomethoxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)

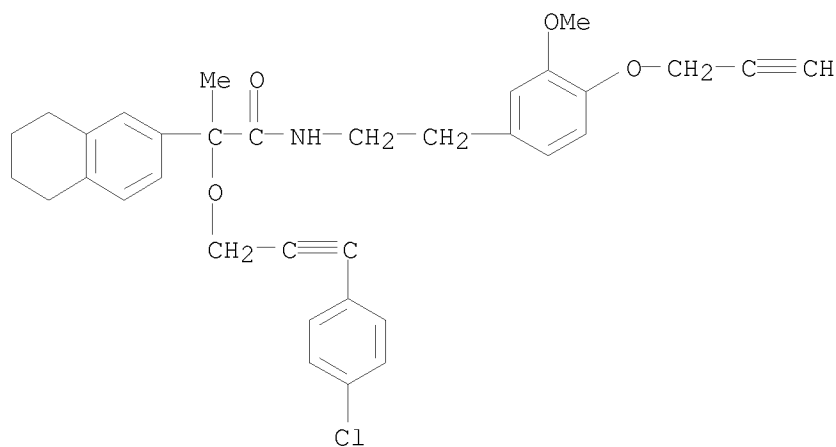
10/513699



RN 1055218-09-1 CAPLUS
CN 2-Naphthaleneacetamide, α -[[3-(4-chlorophenyl)-2-propyn-1-yl]oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055218-10-4 CAPLUS
CN 2-Naphthaleneacetamide, α -[[3-(4-chlorophenyl)-2-propyn-1-yl]oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



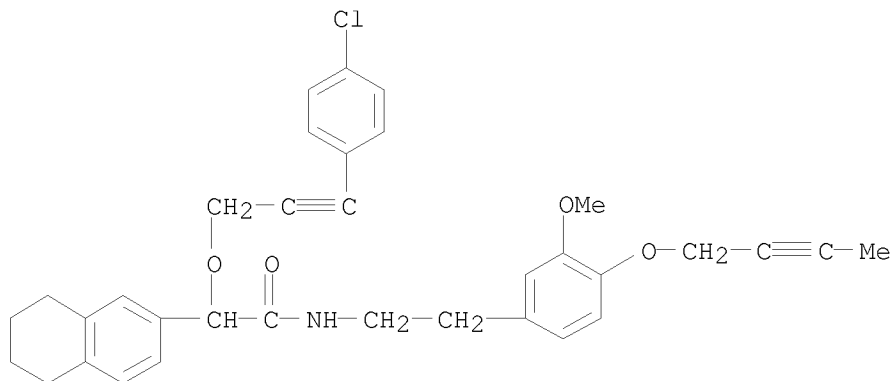
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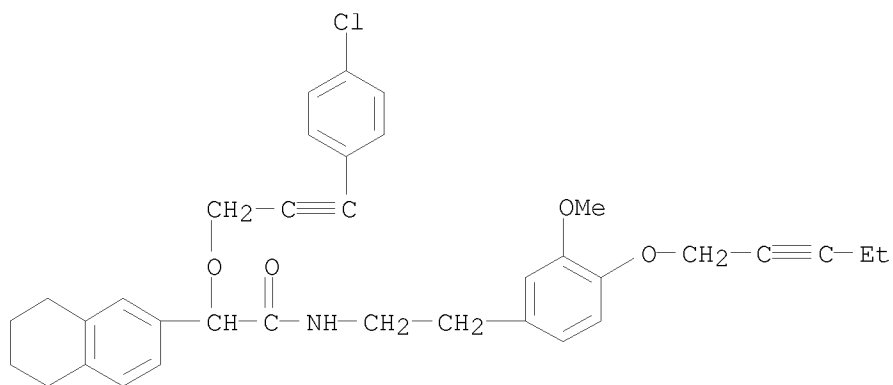
Erich Leese

10/513699

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-
 α -[[3-(4-chlorophenyl)-2-propyn-1-yl]oxy]-5,6,7,8-tetrahydro- (CA
INDEX NAME)

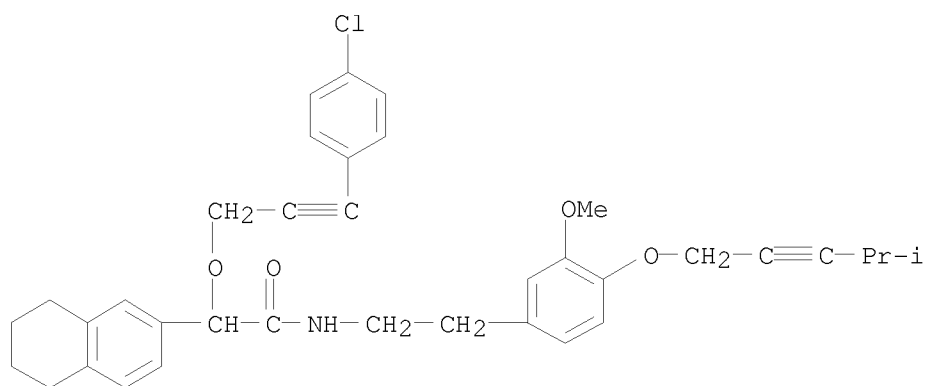


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CN INDEX NAME NOT YET ASSIGNED

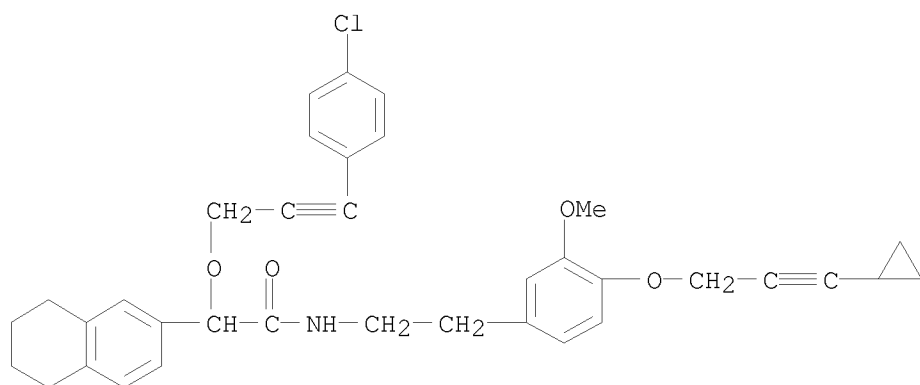


RN 1055218-13-7 CAPLUS
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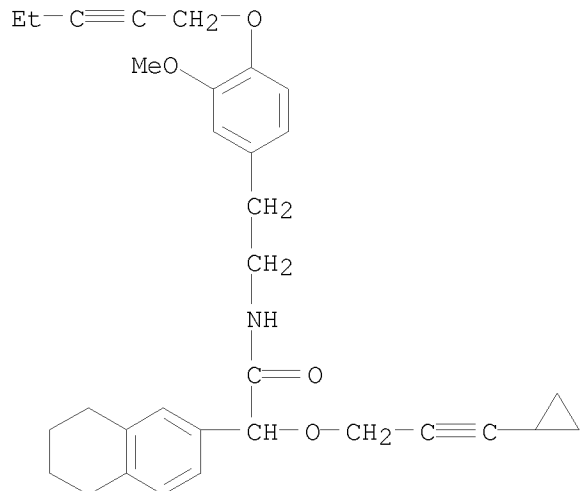


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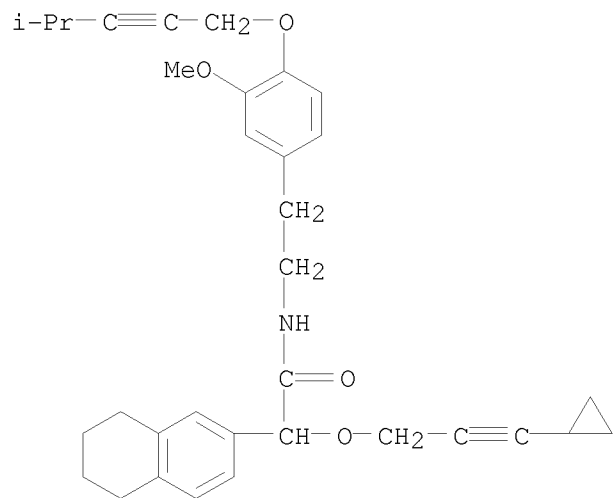


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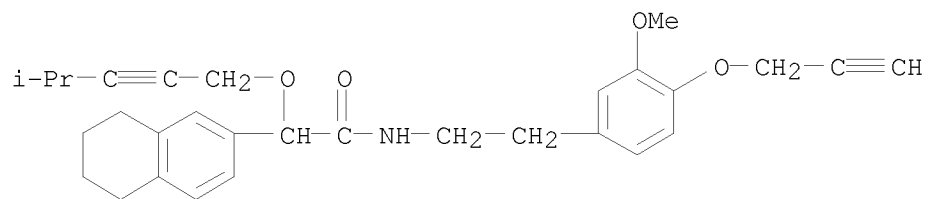
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RN 1055220-19-3 CAPLUS
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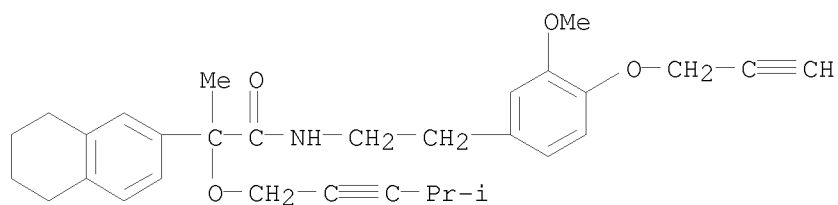
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Erich Leese

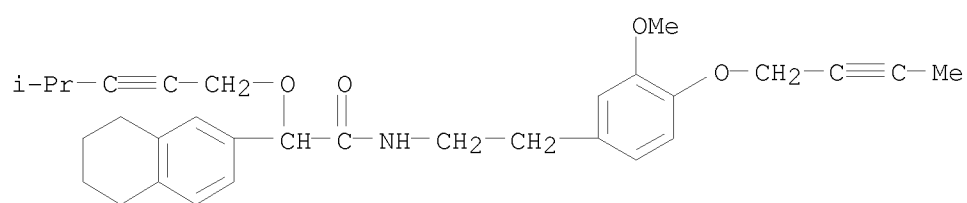
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CN INDEX NAME NOT YET ASSIGNED



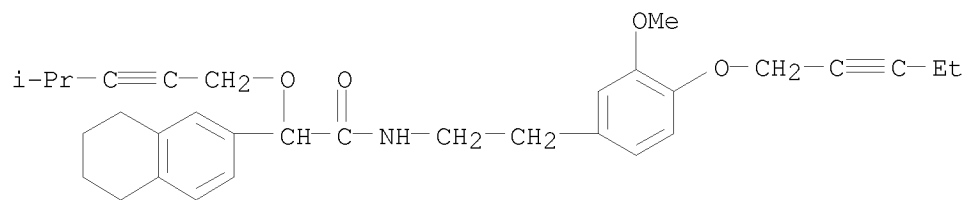
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CN INDEX NAME NOT YET ASSIGNED



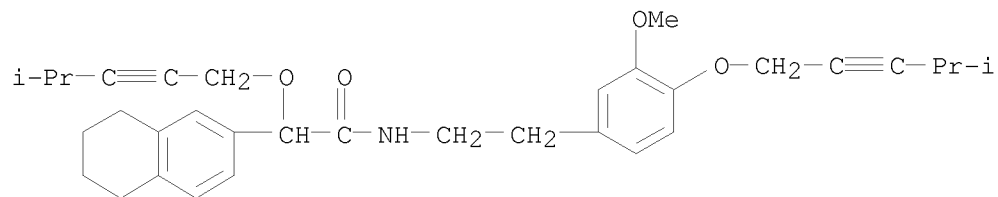
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CN INDEX NAME NOT YET ASSIGNED



RN 1055220-23-9 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



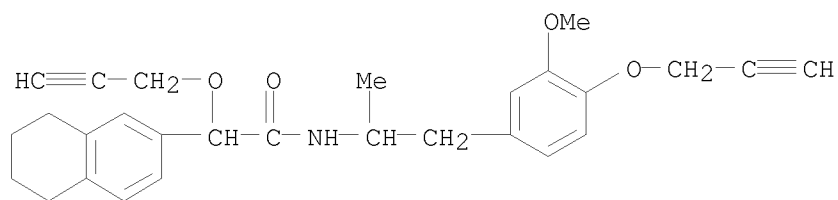
RN 1055221-97-0 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]-1-methylethyl]-α-(2-propyn-1-yloxy)- (CA INDEX NAME)

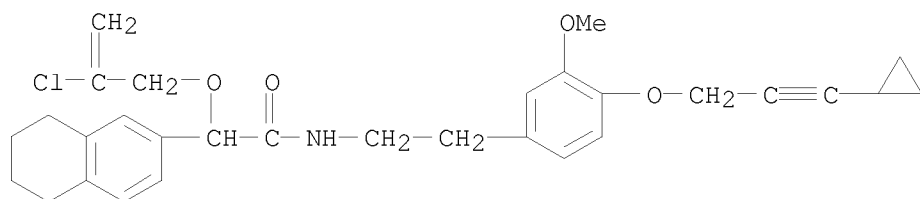
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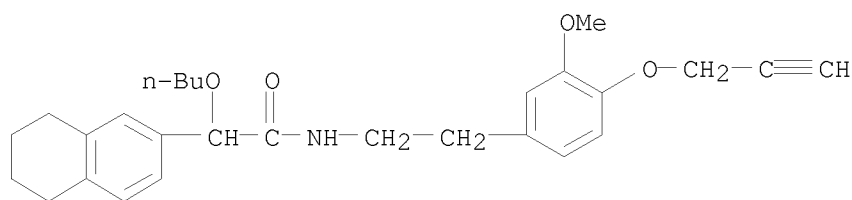
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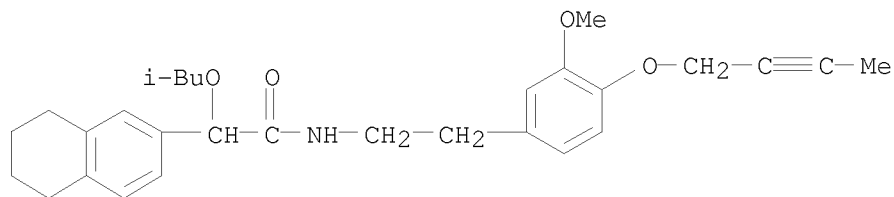
RN 1055222-32-6 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055223-53-4 CAPLUS
CN 2-Naphthaleneacetamide, α -butoxy-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055224-57-1 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-methylpropoxy)- (CA INDEX NAME)



RN 1055224-58-2 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- α -(2-methylpropoxy)- (CA INDEX NAME)

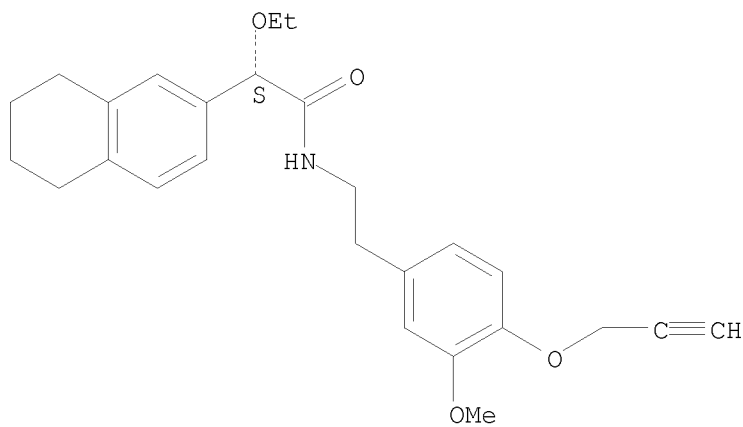
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CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]- α -ethoxy-5,6,7,8-tetrahydro- (CA INDEX NAME)

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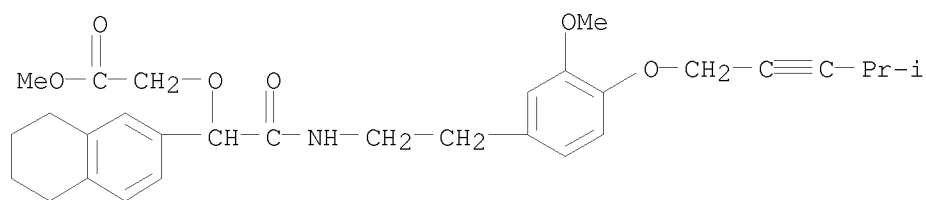
CN 2-Naphthaleneacetamide, α -ethoxy-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-, (α S)- (CA INDEX NAME)

Absolute stereochemistry.

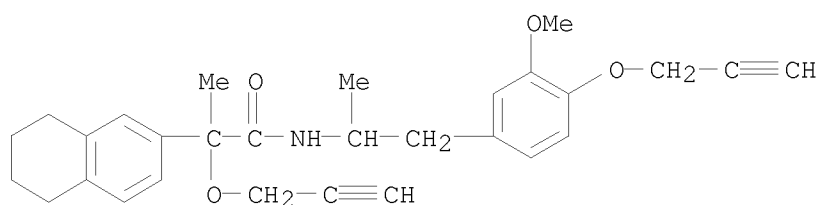


CN INDEX NAME NOT YET ASSIGNED

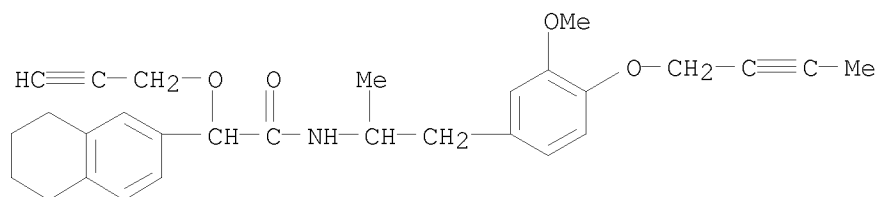
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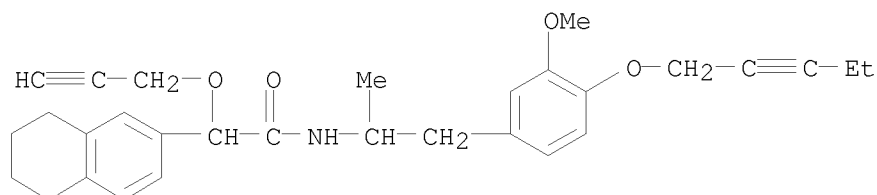
RN 1055228-55-1 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]-1-methylethyl]-alpha-methyl-alpha-(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055228-56-2 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]-1-methylethyl]-5,6,7,8-tetrahydro-alpha-(2-propyn-1-yloxy)- (CA INDEX NAME)

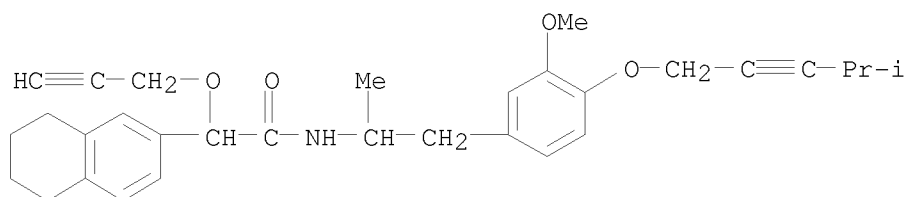


RN 1055228-57-3 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]-1-methylethyl]-alpha-(2-propyn-1-yloxy)- (CA INDEX NAME)



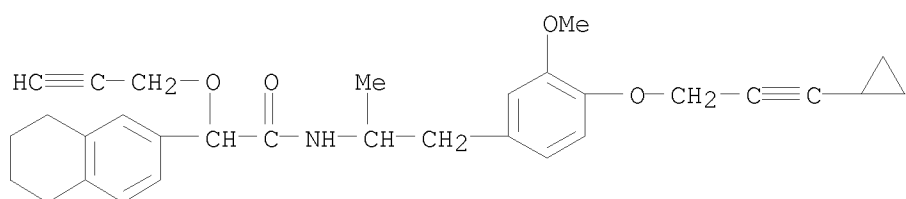
RN 1055228-58-4 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

10/513699



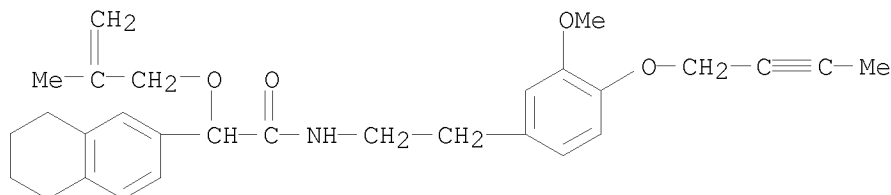
RN 1055228-59-5 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]-1-methylethyl]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



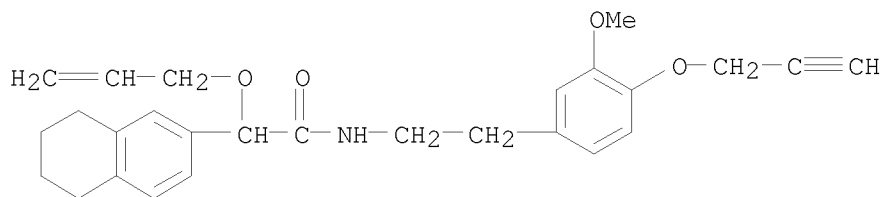
RN 1055229-27-0 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -[(2-methyl-2-propen-1-yl)oxy]- (CA INDEX NAME)



RN 1055230-40-4 CAPLUS

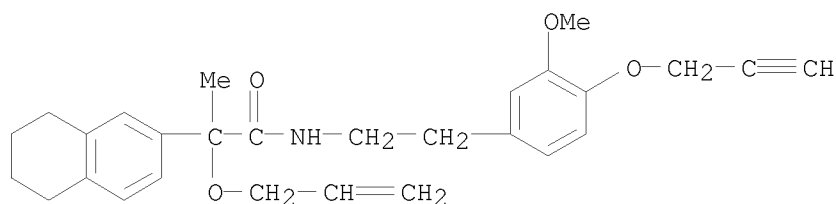
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(2-propen-1-yloxy)- (CA INDEX NAME)



RN 1055230-41-5 CAPLUS

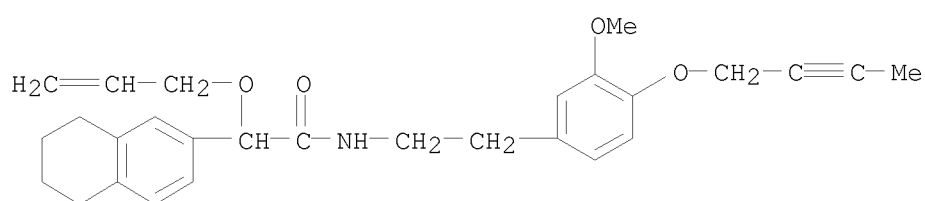
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- α -(2-propen-1-yloxy)- (CA INDEX NAME)

10/513699



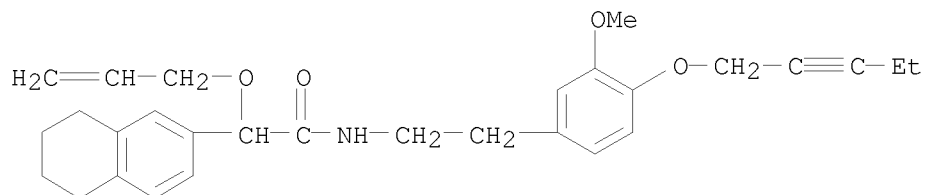
RN 1055230-42-6 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-propen-1-yloxy)- (CA INDEX NAME)



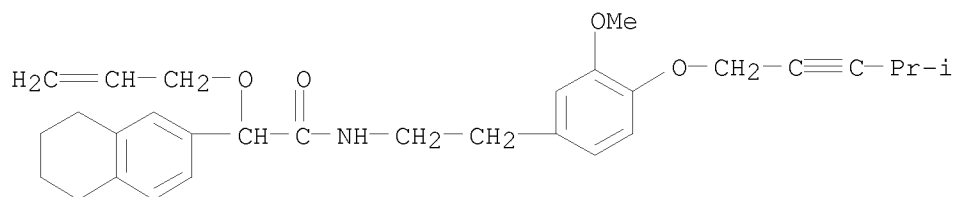
RN 1055230-43-7 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- α -(2-propen-1-yloxy)- (CA INDEX NAME)



RN 1055230-44-8 CAPLUS

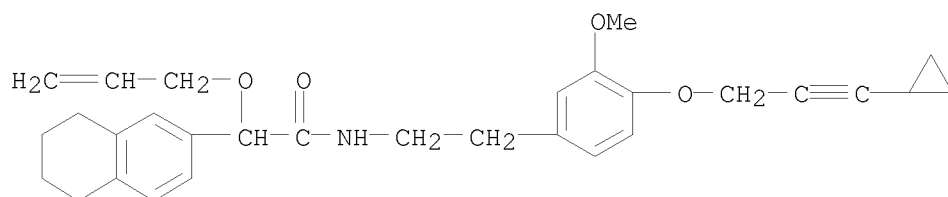
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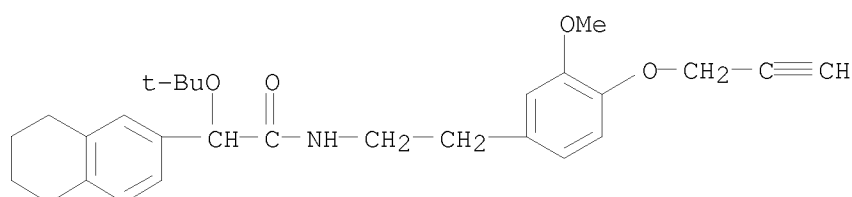
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-propen-1-yloxy)- (CA INDEX NAME)

10/513699



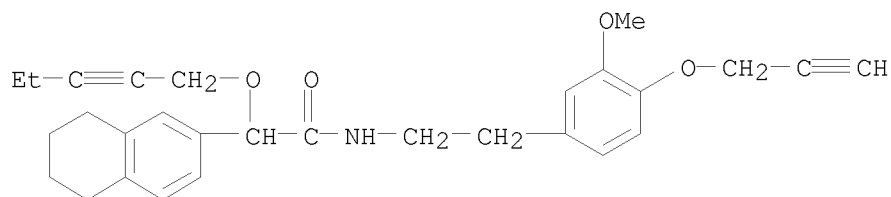
RN 1055230-72-2 CAPLUS

CN 2-Naphthaleneacetamide, α -(1,1-dimethylethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



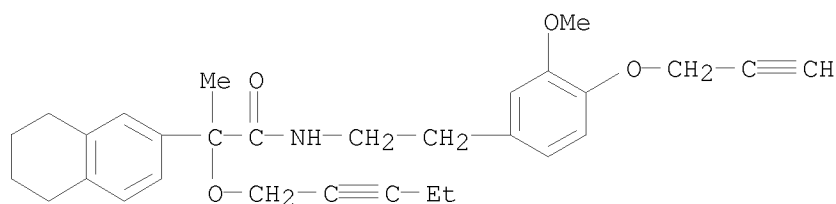
RN 1055230-99-3 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(2-pentyn-1-yloxy)- (CA INDEX NAME)



RN 1055231-00-9 CAPLUS

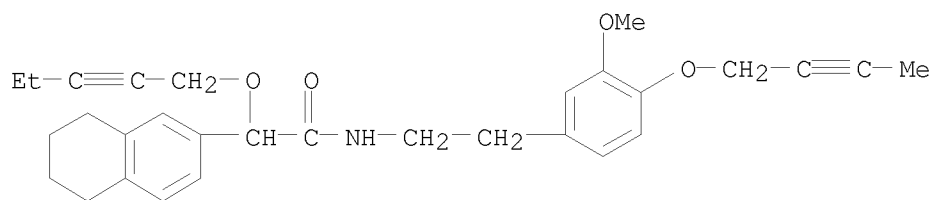
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- α -(2-pentyn-1-yloxy)- (CA INDEX NAME)



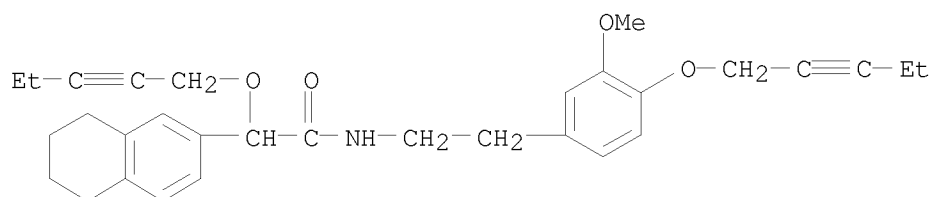
RN 1055231-01-0 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-pentyn-1-yloxy)- (CA INDEX NAME)

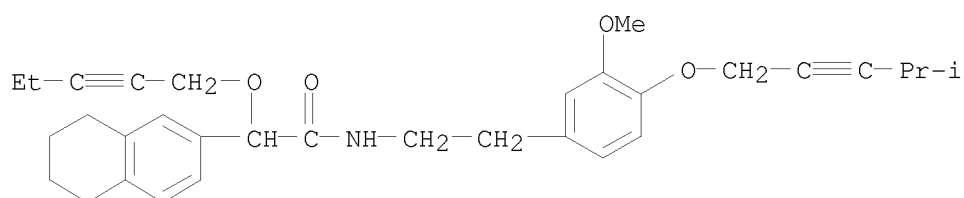
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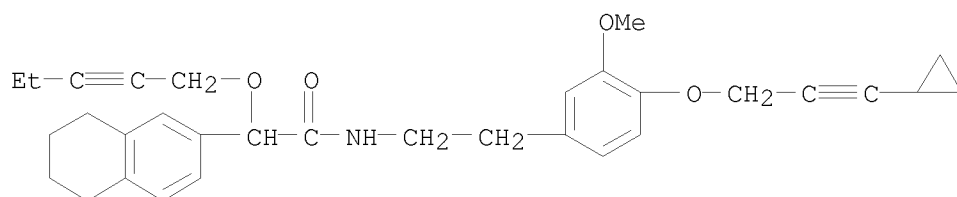
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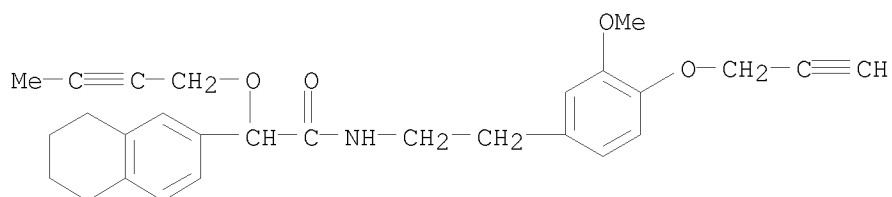


RN 1055231-04-3 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



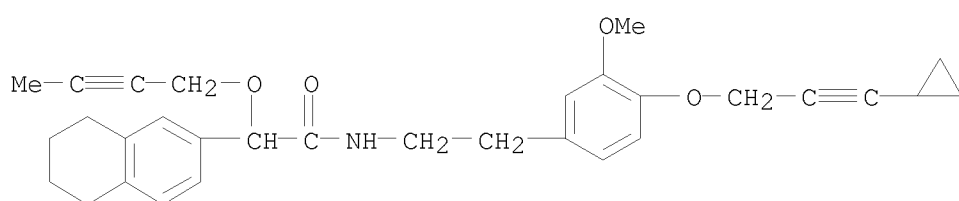
RN 1055232-70-6 CAPLUS
CN 2-Naphthaleneacetamide, α -(2-butyn-1-yloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

10/513699



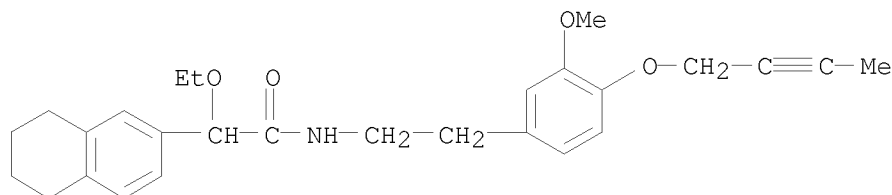
RN 1055232-71-7 CAPLUS

CN 2-Naphthaleneacetamide, α -(2-butyn-1-yloxy)-N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)



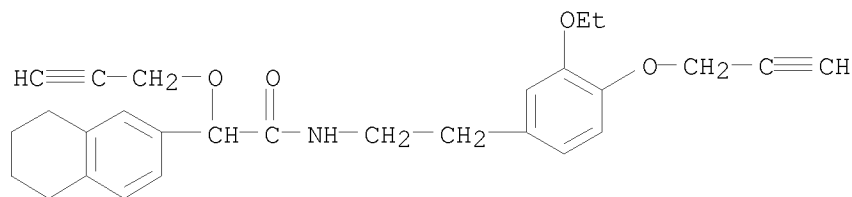
RN 1055236-78-6 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -ethoxy-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 1055238-63-5 CAPLUS

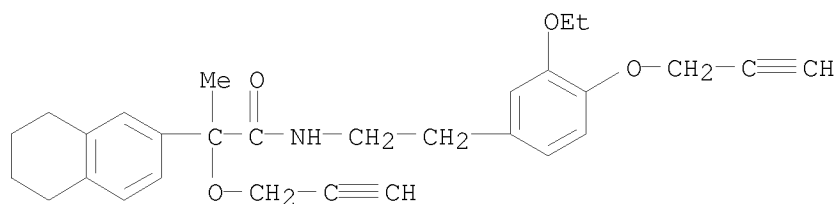
CN 2-Naphthaleneacetamide, N-[2-[3-ethoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055238-64-6 CAPLUS

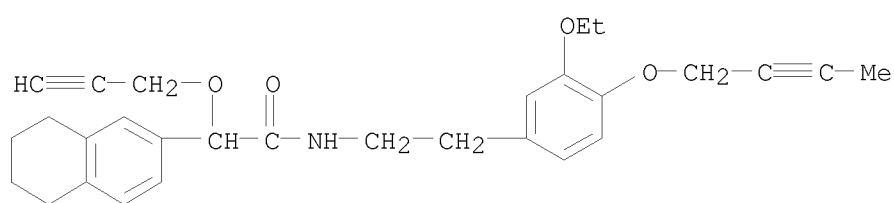
CN 2-Naphthaleneacetamide, N-[2-[3-ethoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-5,6,7,8-tetrahydro- α -methyl- α -(2-propyn-1-yloxy)- (CA INDEX NAME)

10/513699



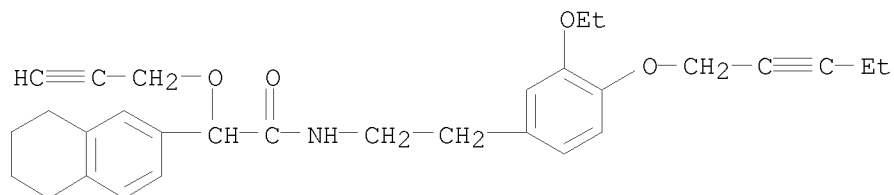
RN 1055238-65-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-ethoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



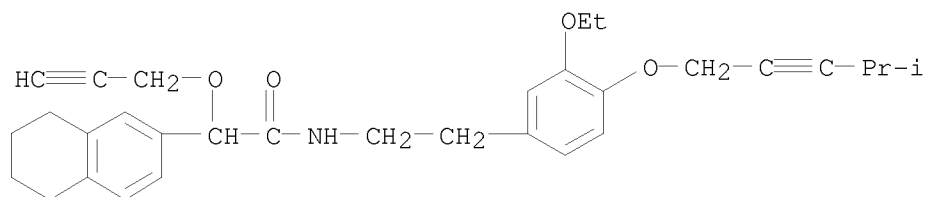
RN 1055238-66-8 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[3-ethoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055238-67-9 CAPLUS

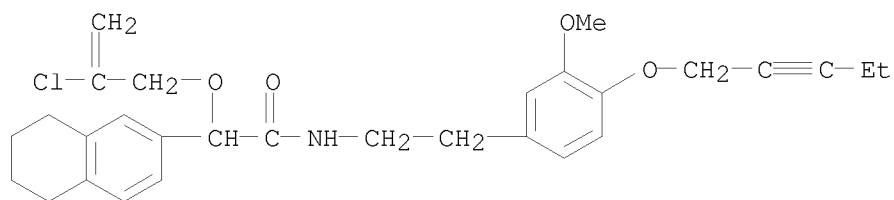
CN INDEX NAME NOT YET ASSIGNED



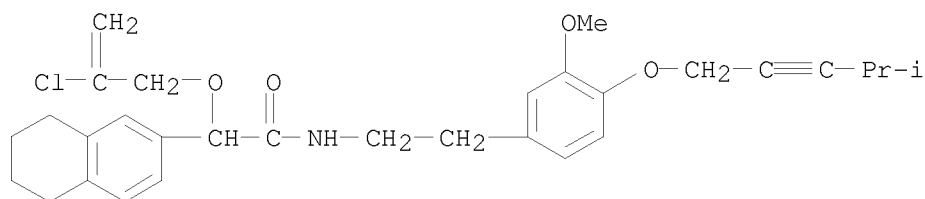
RN 1055240-59-9 CAPLUS

CN INDEX NAME NOT YET ASSIGNED

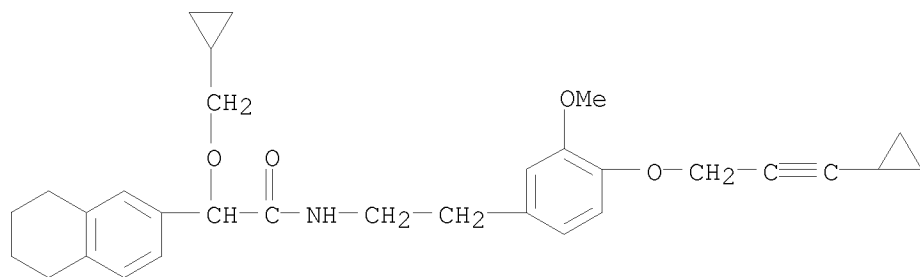
10/513699



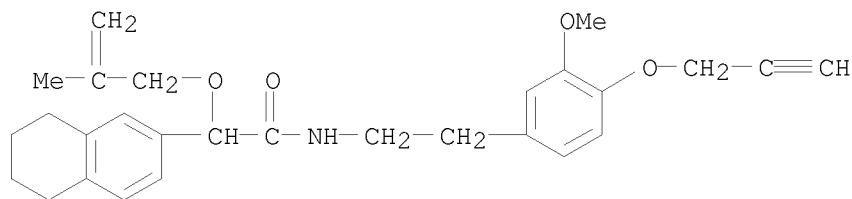
RN 1055240-60-2 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055242-23-3 CAPLUS
CN 2-Naphthaleneacetamide, α -(cyclopropylmethoxy)-N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)



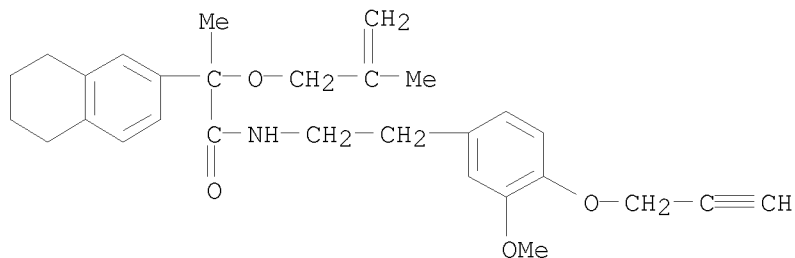
RN 1055243-85-0 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(2-methyl-2-propen-1-yl)oxy- (CA INDEX NAME)



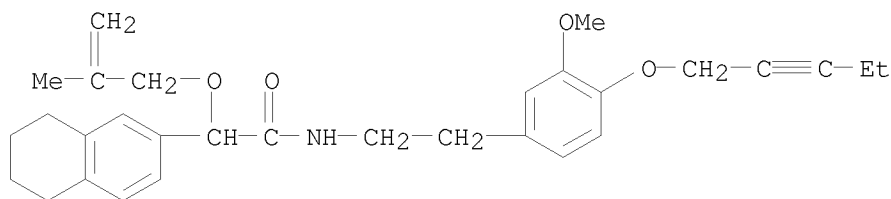
RN 1055243-86-1 CAPLUS

10/513699

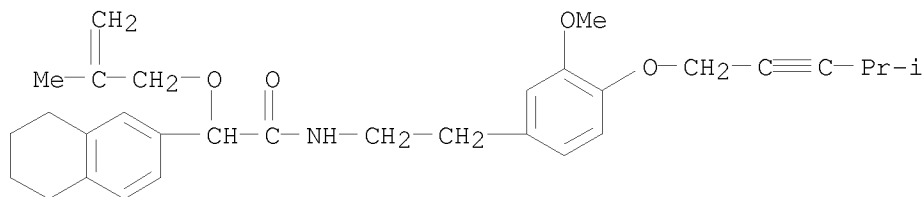
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- α -(2-methyl-2-propen-1-yl)oxy]-
(CA INDEX NAME)



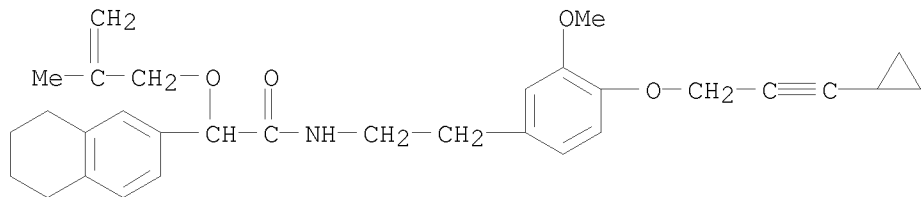
RN 1055243-87-2 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055243-88-3 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



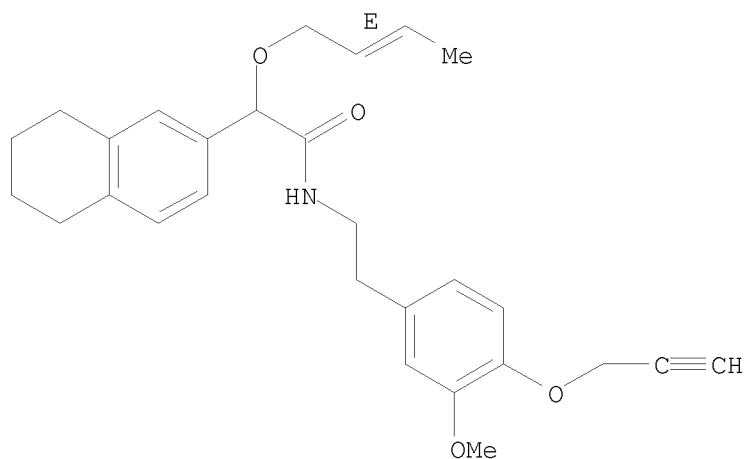
RN 1055243-89-4 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-methyl-2-propen-1-yl)oxy]- (CA INDEX NAME)



10/513699

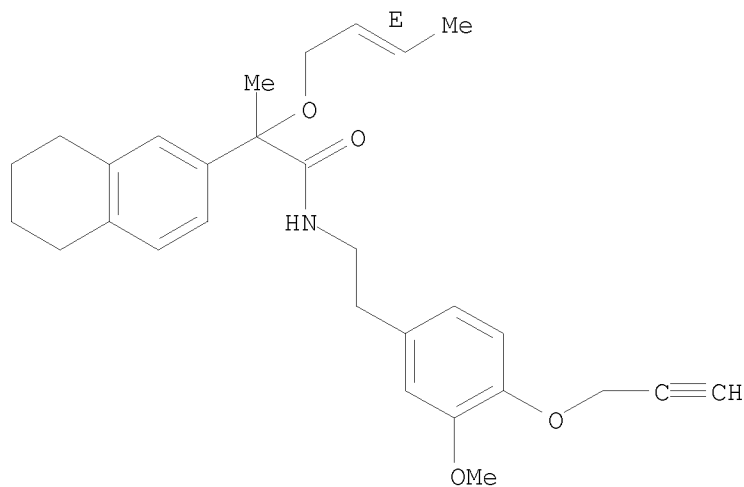
RN 1055244-11-5 CAPLUS
CN 2-Naphthaleneacetamide, α -[(2E)-2-buten-1-yloxy]-5,6,7,8-tetrahydro-
N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

Double bond geometry as shown.



RN 1055244-12-6 CAPLUS
CN 2-Naphthaleneacetamide, α -[(2E)-2-buten-1-yloxy]-5,6,7,8-tetrahydro-
N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA
INDEX NAME)

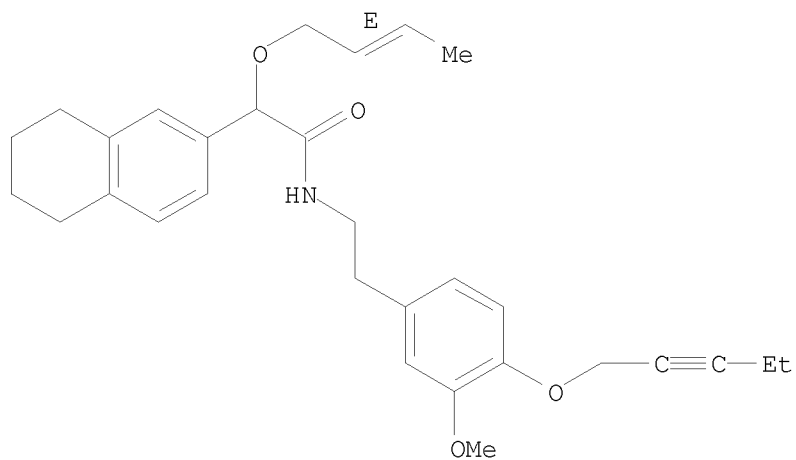
Double bond geometry as shown.



RN 1055244-13-7 CAPLUS
CN 2-Naphthaleneacetamide, α -[(2E)-2-buten-1-yloxy]-5,6,7,8-tetrahydro-
N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

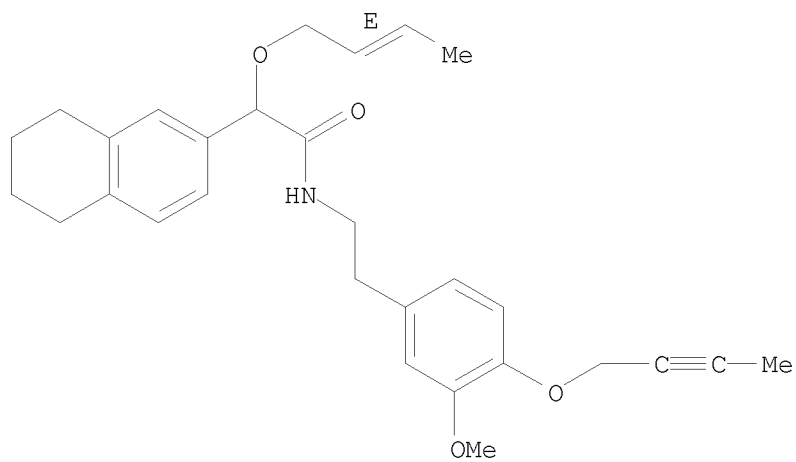
Double bond geometry as shown.

10/513699



RN 1055244-14-8 CAPLUS
CN 2-Naphthaleneacetamide, α -[(2E)-2-buten-1-yloxy]-N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)

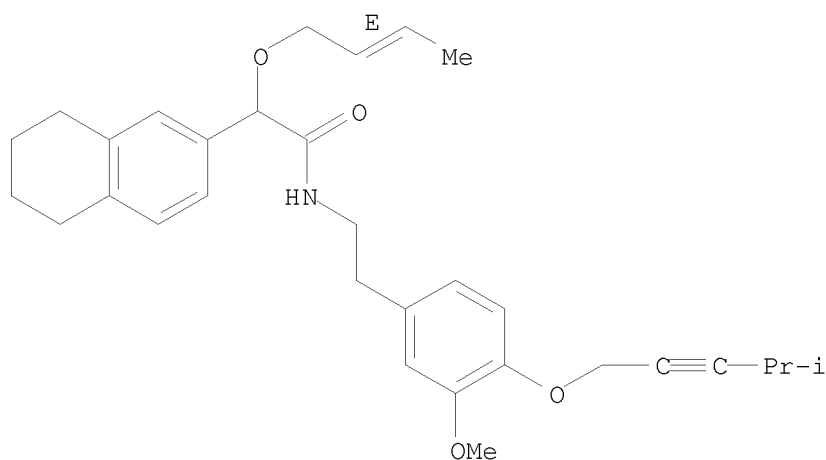
Double bond geometry as shown.



RN 1055244-15-9 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

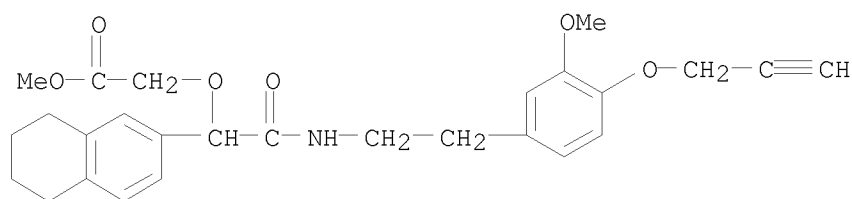
Double bond geometry as shown.

10/513699



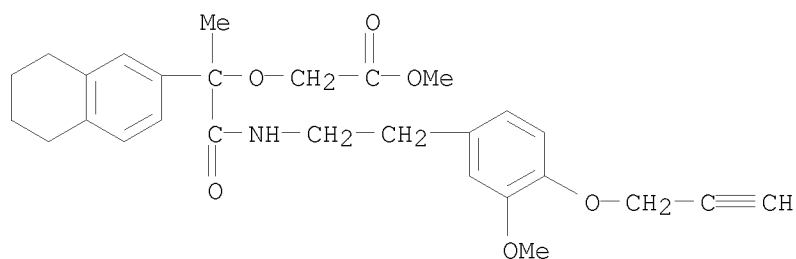
RN 1055247-12-5 CAPLUS

CN Acetic acid, 2-[2-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-2-oxo-1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethoxy]-, methyl ester (CA INDEX NAME)



RN 1055247-15-8 CAPLUS

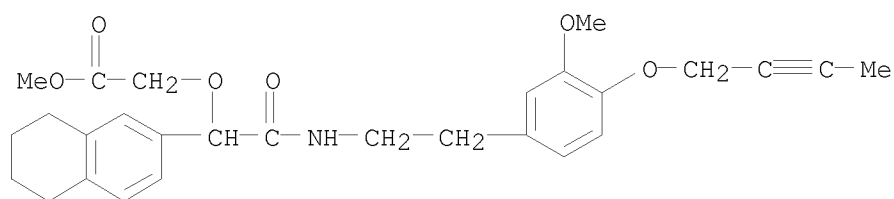
CN Acetic acid, 2-[2-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-1-methyl-2-oxo-1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethoxy]-, methyl ester (CA INDEX NAME)



RN 1055247-16-9 CAPLUS

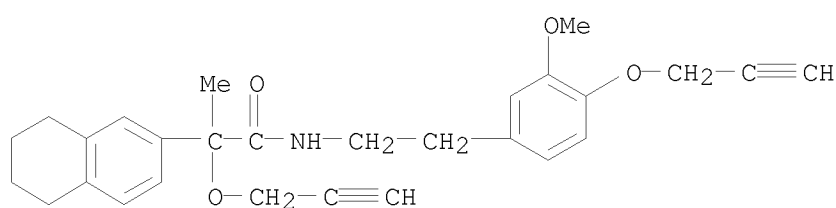
CN Acetic acid, 2-[2-[[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]amino]-2-oxo-1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethoxy]-, methyl ester (CA INDEX NAME)

10/513699



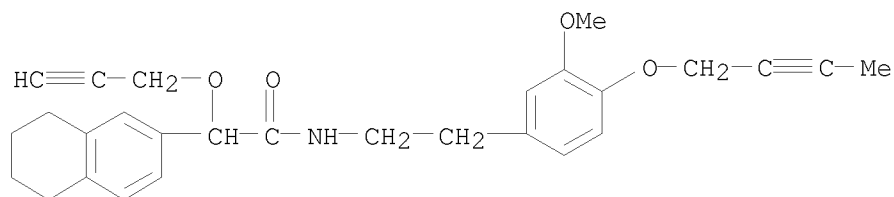
RN 1055249-28-9 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-(2-propyn-1-yloxy)- (CA INDEX NAME)



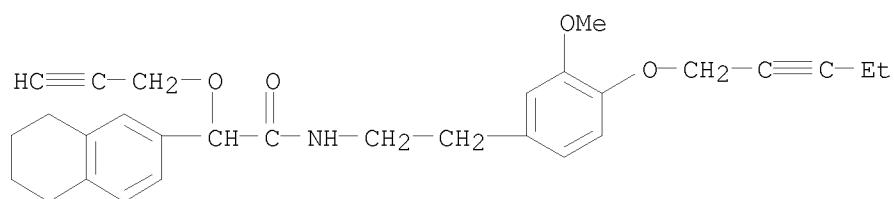
RN 1055249-29-0 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055249-30-3 CAPLUS

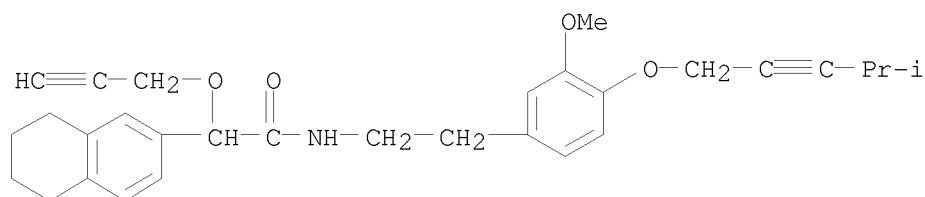
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]-α-(2-propyn-1-yloxy)- (CA INDEX NAME)



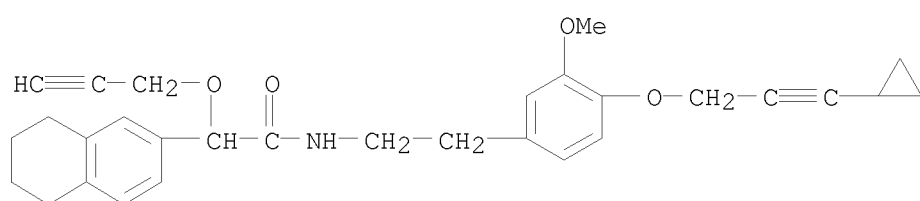
RN 1055249-33-6 CAPLUS

CN INDEX NAME NOT YET ASSIGNED

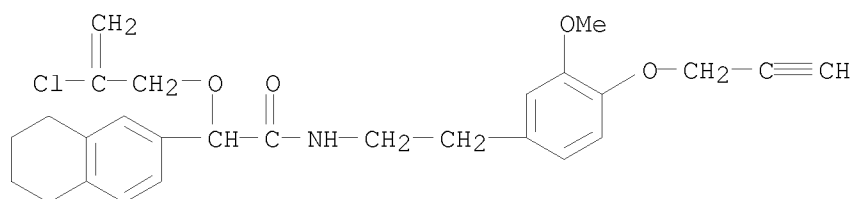
10/513699



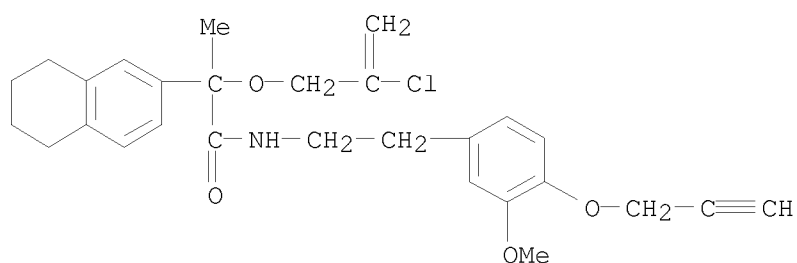
RN 1055249-34-7 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055250-39-9 CAPLUS
CN 2-Naphthaleneacetamide, α -[(2-chloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



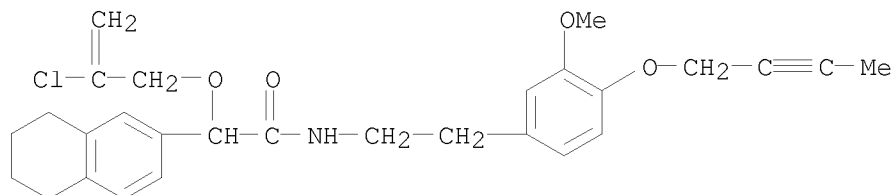
RN 1055250-41-3 CAPLUS
CN 2-Naphthaleneacetamide, α -[(2-chloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



10/513699

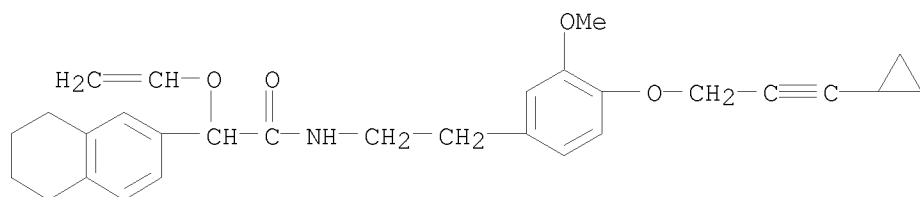
RN 1055250-45-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-
 α -(2-chloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro- (CA INDEX NAME)



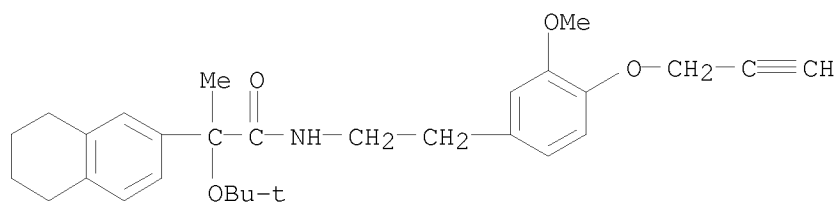
RN 1055253-12-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]- α -(ethenyloxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)



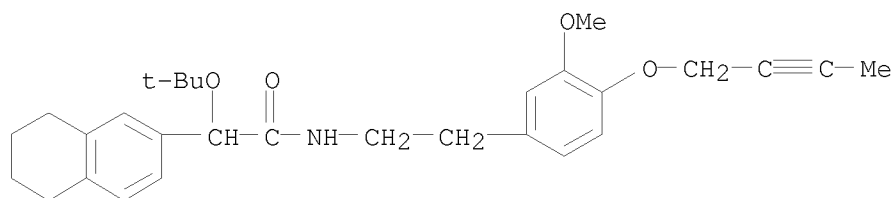
RN 1055254-60-8 CAPLUS

CN 2-Naphthaleneacetamide, α -(1,1-dimethylethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



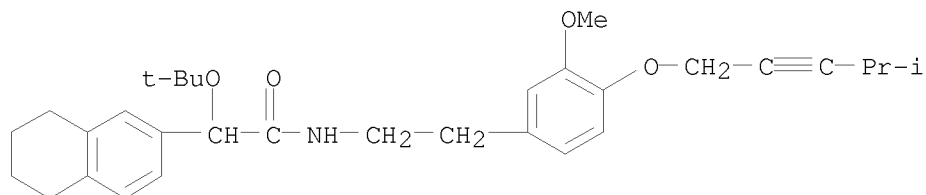
RN 1055254-61-9 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-
 α -(1,1-dimethylethoxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)

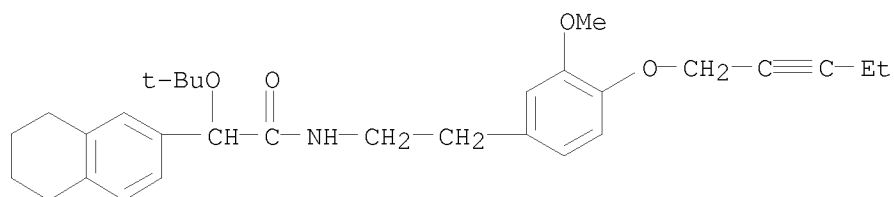


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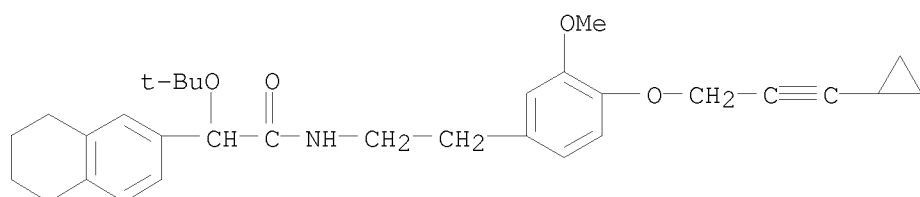
RN 1055254-62-0 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



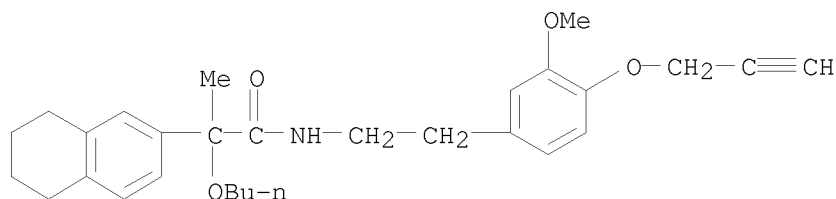
RN 1055254-63-1 CAPLUS
CN 2-Naphthaleneacetamide, α -(1,1-dimethylethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055254-64-2 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]- α -(1,1-dimethylethoxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)

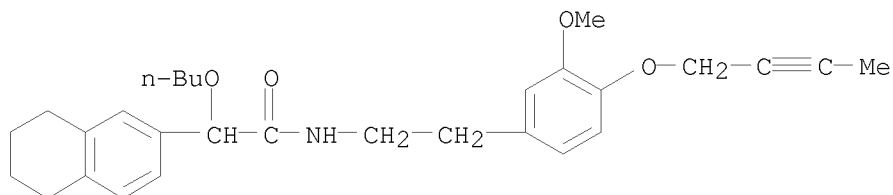


RN 1055255-69-0 CAPLUS
CN 2-Naphthaleneacetamide, α -butoxy-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)

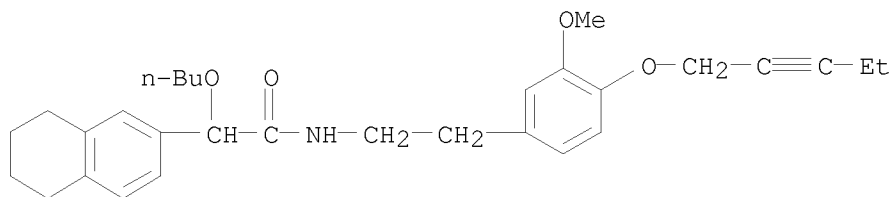


10/513699

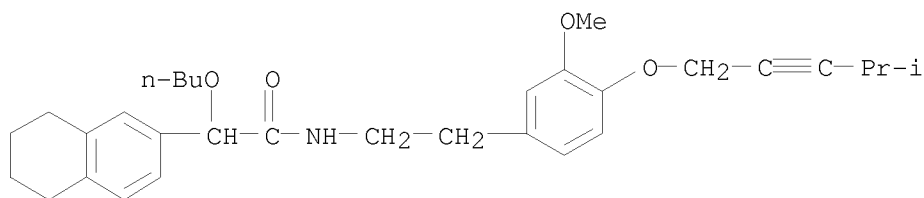
RN 1055255-70-3 CAPLUS
CN 2-Naphthaleneacetamide, α -butoxy-N-[2-[4-(2-butyne-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)



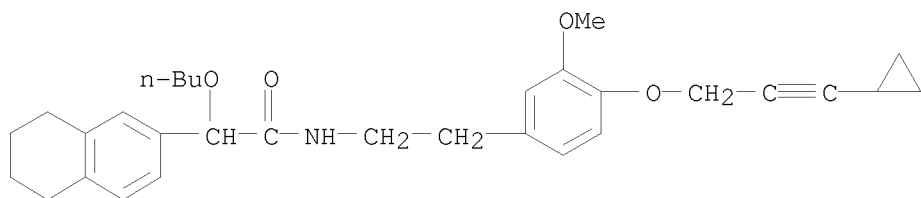
RN 1055255-71-4 CAPLUS
CN 2-Naphthaleneacetamide, α -butoxy-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyne-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055255-72-5 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055255-73-6 CAPLUS
CN 2-Naphthaleneacetamide, α -butoxy-N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)

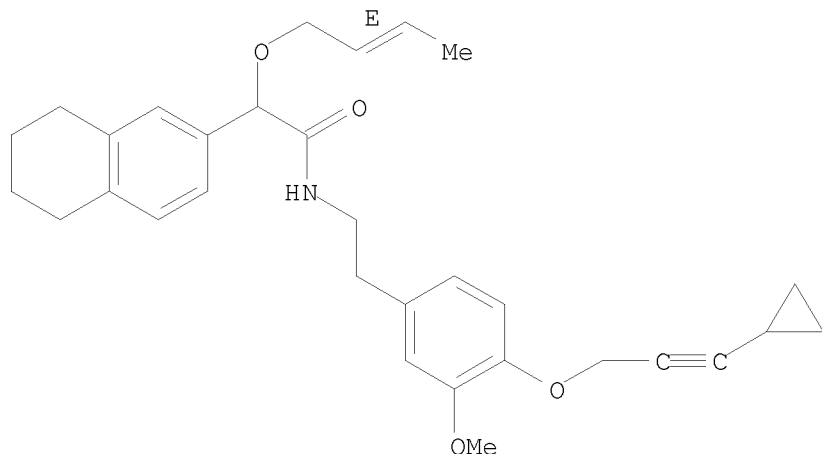


RN 1055256-73-9 CAPLUS
CN 2-Naphthaleneacetamide, α -[(2E)-2-buten-1-yloxy]-N-[2-[4-[(3-

10/513699

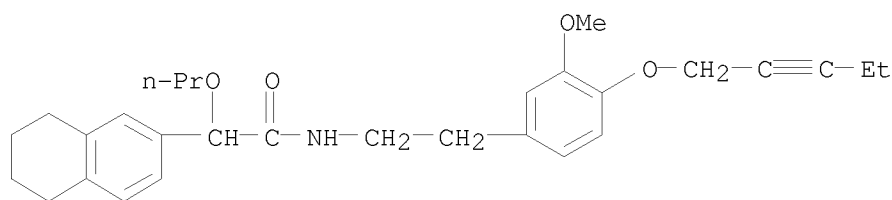
cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-
(CA INDEX NAME)

Double bond geometry as shown.



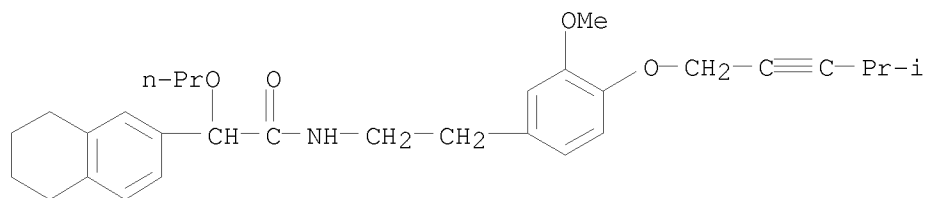
RN 1055258-31-5 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]-α-propoxy- (CA INDEX NAME)



RN 1055258-32-6 CAPLUS

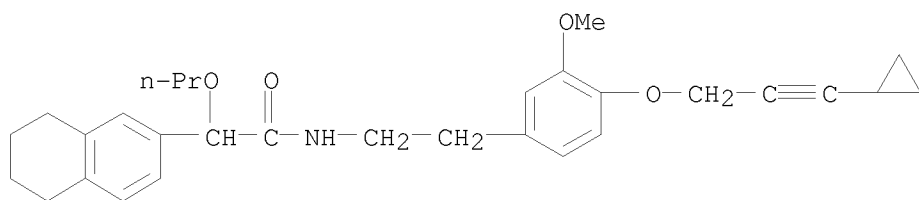
CN INDEX NAME NOT YET ASSIGNED



RN 1055258-33-7 CAPLUS

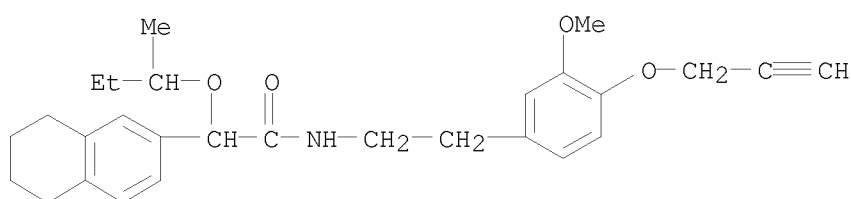
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-propoxy- (CA INDEX NAME)

10/513699



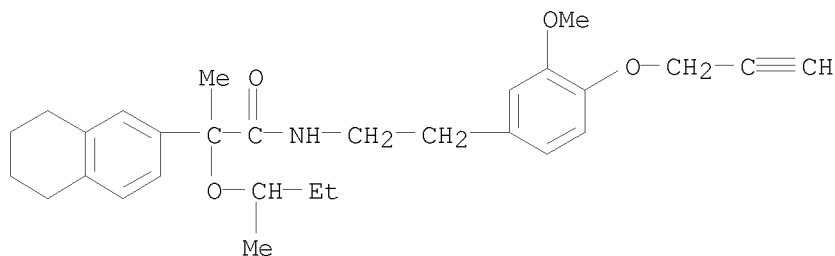
RN 1055261-20-5 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(1-methylpropoxy)- (CA INDEX NAME)



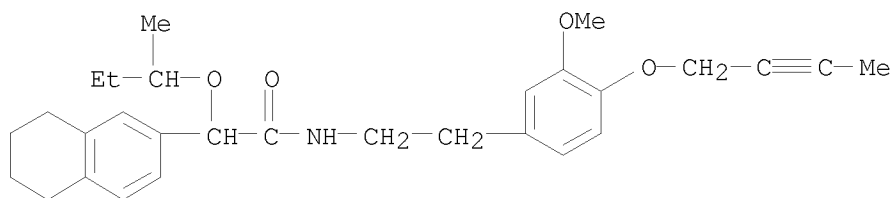
RN 1055261-21-6 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-(1-methylpropoxy)- (CA INDEX NAME)



RN 1055261-22-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(1-methylpropoxy)- (CA INDEX NAME)

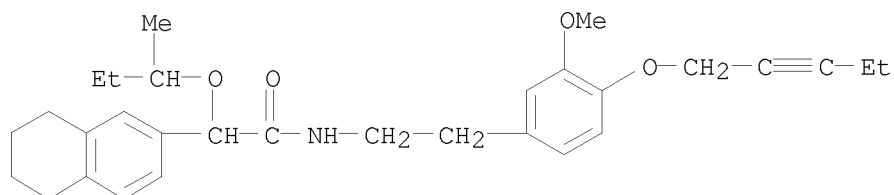


RN 1055261-23-8 CAPLUS

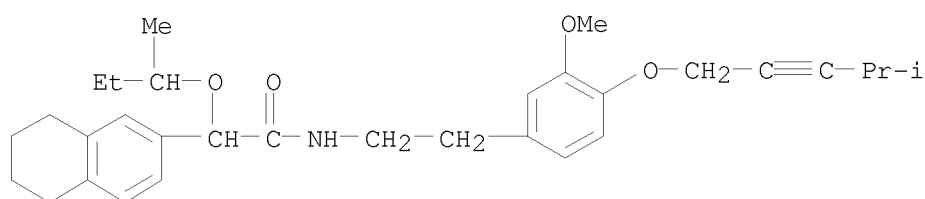
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]-α-methyl-α-(1-methylpropoxy)- (CA INDEX NAME)

10/513699

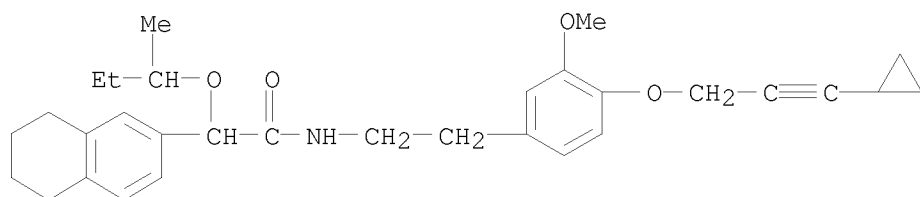
yloxy)phenyl]ethyl]- α -(1-methylpropoxy)- (CA INDEX NAME)



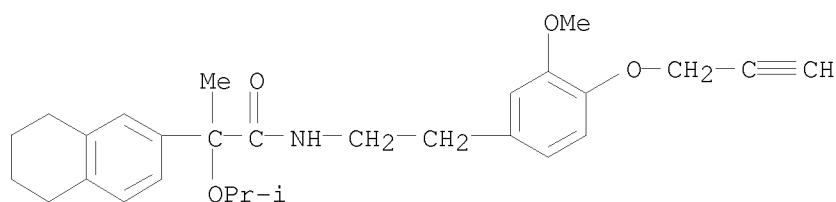
RN 1055261-24-9 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055261-25-0 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(1-methylpropoxy)- (CA INDEX NAME)



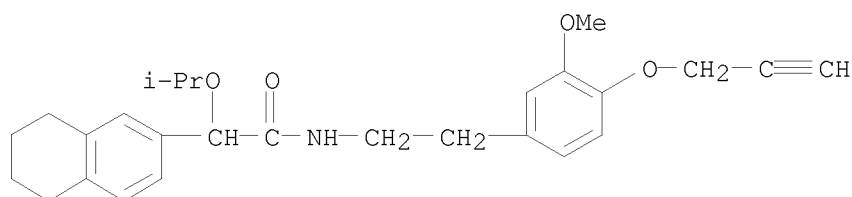
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CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- α -(1-methylethoxy)- (CA INDEX NAME)



RN 1055262-70-8 CAPLUS
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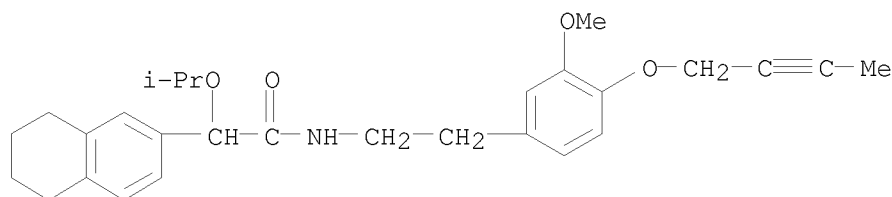
10/513699

yloxy)phenyl]ethyl]- α -(1-methylethoxy)- (CA INDEX NAME)



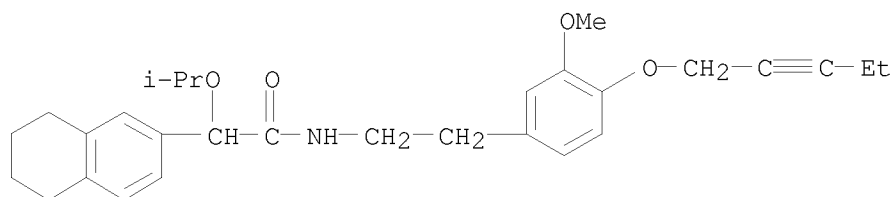
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CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(1-methylethoxy)- (CA INDEX NAME)



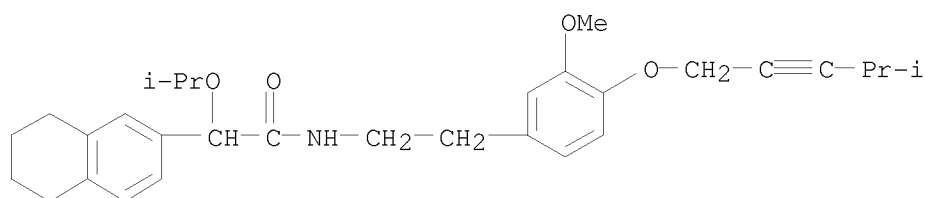
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CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- α -(1-methylethoxy)- (CA INDEX NAME)



RN 1055262-76-4 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



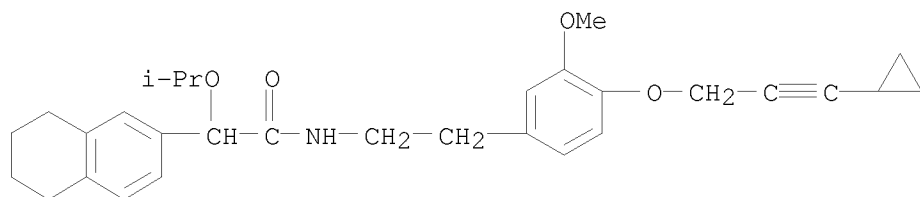
RN 1055262-78-6 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(1-methylethoxy)- (CA INDEX NAME)

<12/04/2007>

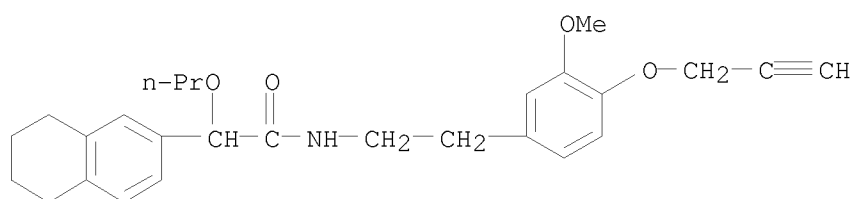
Erich Leese

10/513699



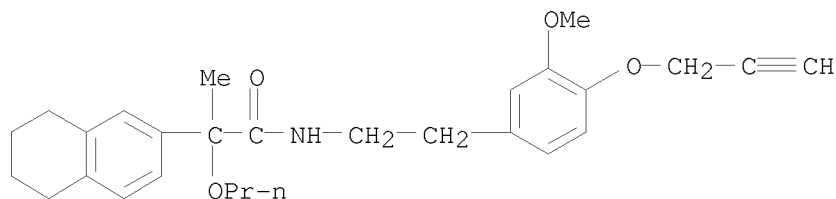
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CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -propoxy- (CA INDEX NAME)



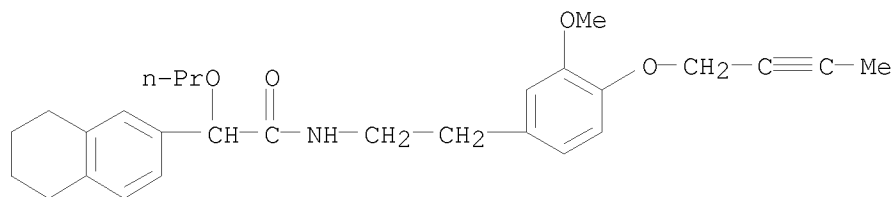
RN 1055264-43-1 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- α -propoxy- (CA INDEX NAME)



RN 1055264-45-3 CAPLUS

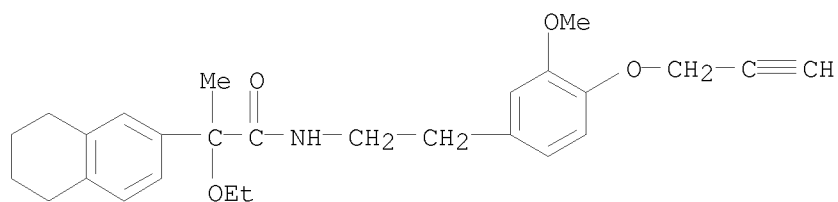
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -propoxy- (CA INDEX NAME)



RN 1055267-02-1 CAPLUS

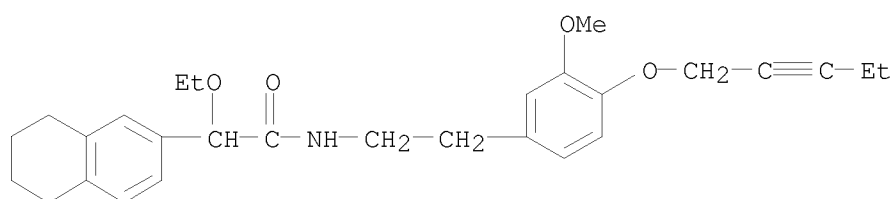
CN 2-Naphthaleneacetamide, α -ethoxy-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)

10/513699



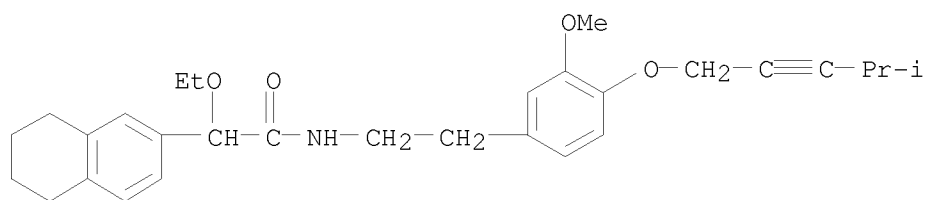
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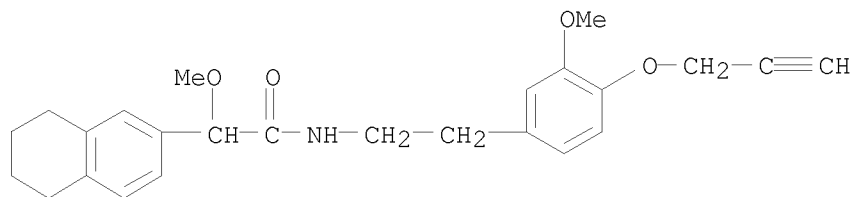
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CN INDEX NAME NOT YET ASSIGNED



RN 1055270-60-4 CAPLUS

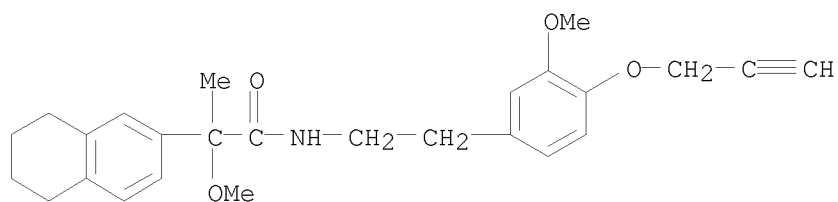
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -methoxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055270-61-5 CAPLUS

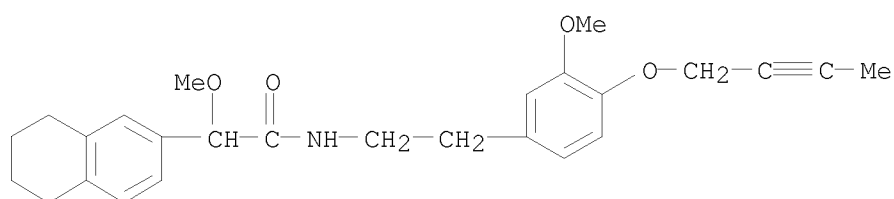
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -methoxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)

10/513699



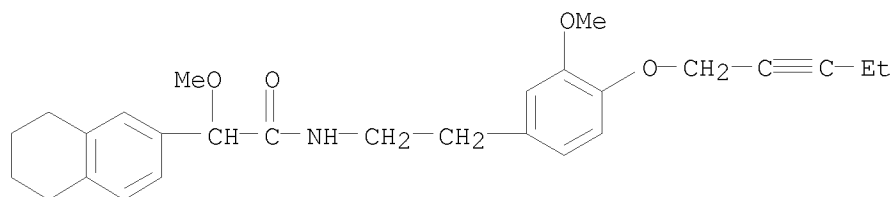
RN 1055270-62-6 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -methoxy- (CA INDEX NAME)



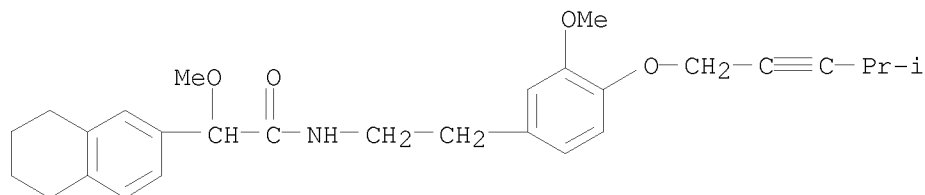
RN 1055270-63-7 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -methoxy-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055270-64-8 CAPLUS

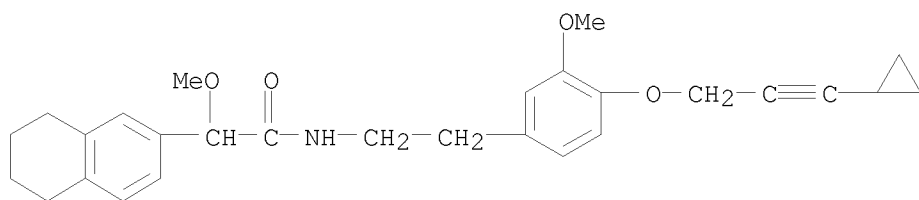
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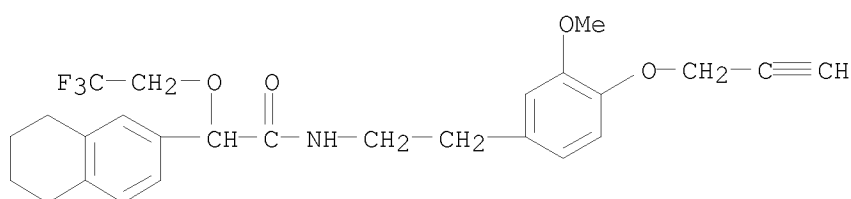
RN 1055270-65-9 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -methoxy- (CA INDEX NAME)

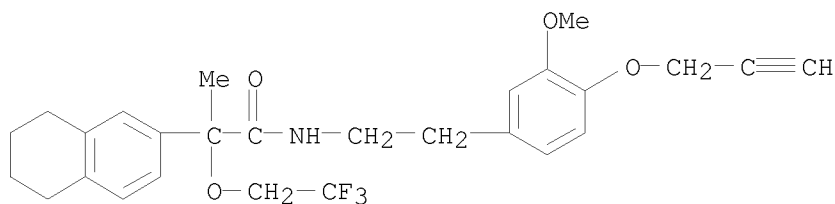
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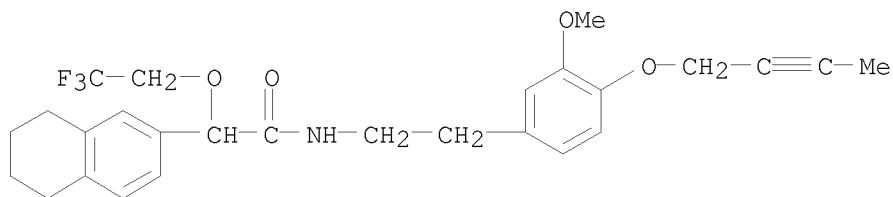
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CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(2,2,2-trifluoroethoxy)- (CA INDEX NAME)



RN 1055271-93-6 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-(2,2,2-trifluoroethoxy)- (CA INDEX NAME)

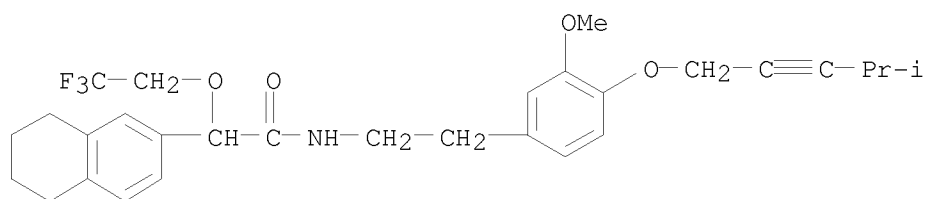


RN 1055271-94-7 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(2,2,2-trifluoroethoxy)- (CA INDEX NAME)



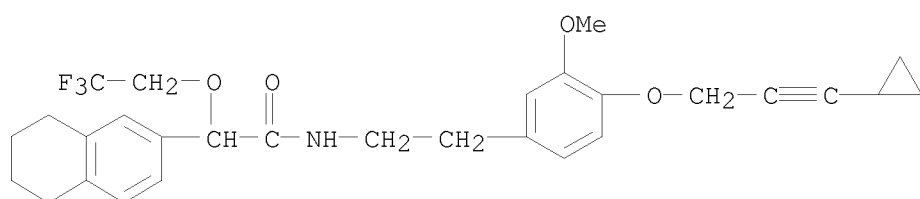
RN 1055271-95-8 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

10/513699



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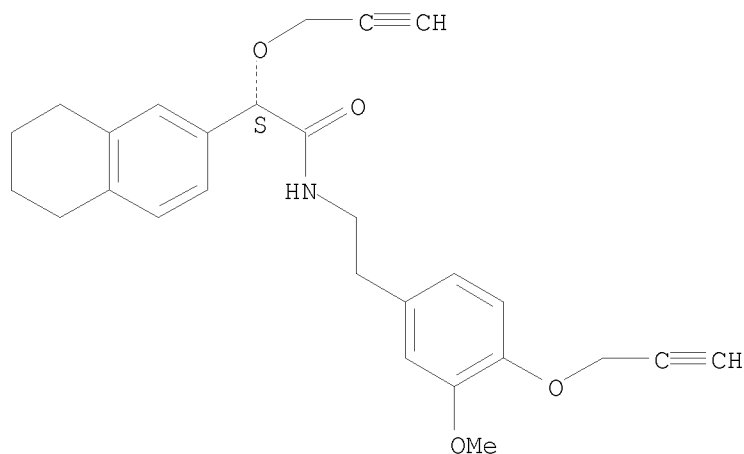
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2,2,2-trifluoroethoxy)- (CA INDEX NAME)



RN 1055272-25-7 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(2-propyn-1-yloxy)-, (α S)- (CA INDEX NAME)

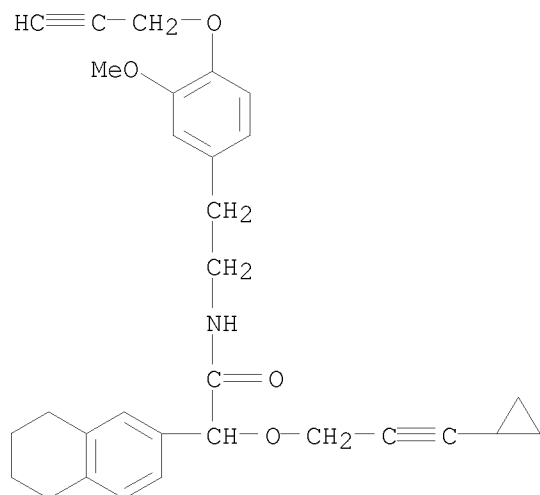
Absolute stereochemistry.



RN 1055273-51-2 CAPLUS

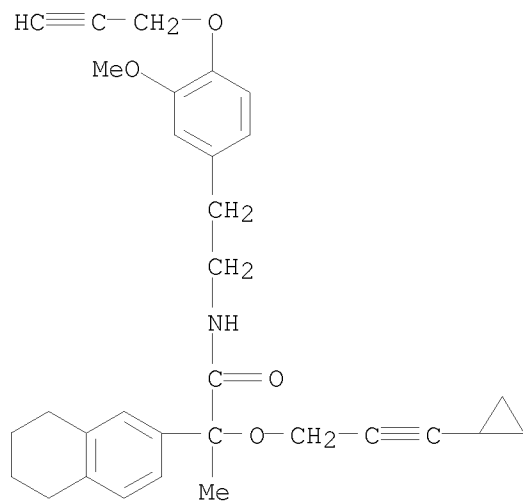
CN 2-Naphthaleneacetamide, α -[(3-cyclopropyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

10/513699



RN 1055273-52-3 CAPLUS

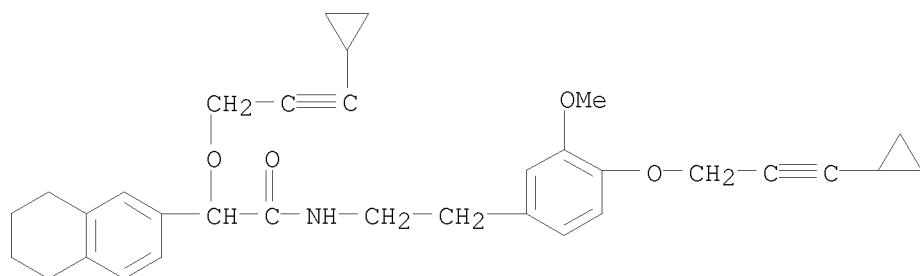
CN 2-Naphthaleneacetamide, α -[(3-cyclopropyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



RN 1055273-53-4 CAPLUS

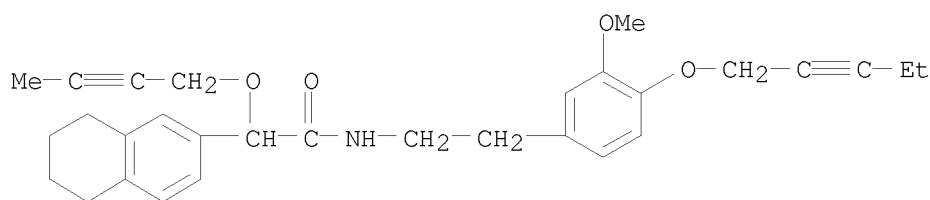
CN 2-Naphthaleneacetamide, α -[(3-cyclopropyl-2-propyn-1-yl)oxy]-N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)

10/513699



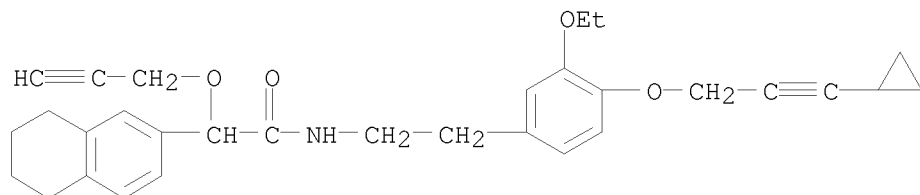
RN 1055274-16-2 CAPLUS

CN 2-Naphthaleneacetamide, α -(2-butyn-1-yloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055275-96-1 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-ethoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



OS.CITING REF COUNT: 1 THERE ARE 1 CAPLUS RECORDS THAT CITE THIS RECORD (1 CITINGS)

REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

10/513699

L5 ANSWER 12 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2002:793595 CAPLUS

DOCUMENT NUMBER: 137:310703

TITLE: Preparation of novel N-propargyloxyphenethyl
thioacetamides as agrochemical fungicides

INVENTOR(S): Kunz, Walter; Lamberth, Clemens; Cederbaum, Fredrik;
Zeller, Martin

PATENT ASSIGNEE(S): Syngenta Participations A.-G., Switz.

SOURCE: PCT Int. Appl., 61 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

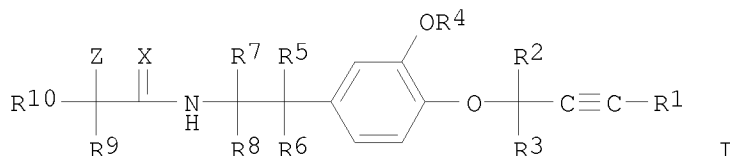
FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2002081437	A2	20021017	WO 2002-EP3623	20020402 <--
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZM, ZW			
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CA 2443131	A1	20021017	CA 2002-2443131	20020402 <--
CA 2443131	C	20090804		
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CN 1514826	A	20040721	CN 2002-807705	20020402 <--
JP 2004526751	T	20040902	JP 2002-579425	20020402 <--
JP 4080891	B2	20080423		
US 20040127739	A1	20040701	US 2003-472577	20030923 <--
US 7105545	B2	20060912		
IN 2003CN01555	A	20051125	IN 2003-CN1555	20031001
MX 2003009091	A	20040212	MX 2003-9091	20031003 <--
PRIORITY APPLN. INFO.:			GB 2001-8339	A 20010403
			WO 2002-EP3623	W 20020402

OTHER SOURCE(S): MARPAT 137:310703

GI



AB The title compds. [I; R1 = H, alkyl, cycloalkyl, (un)substituted aryl; R2,

<12/04/2007>

Erich Leese

R3 = H, alkyl; R4 = alkyl, alkenyl, alkynyl; R5-R8 = H, alkyl; R9 = H, alkyl, alkenyl, alkynyl; R10 = (un)substituted aryl, heteroaryl; Z = OH, (un)substituted aryloxy, (un)substituted alkoxy, etc.; X = S] which possess useful plant protecting properties and may advantageously be employed in agricultural practice for controlling or preventing the infestation of plants by phytopathogenic microorganisms, especially fungi, were prepared. Thus, reacting I [R1-R3 = H; R4 = Me; R5-R9 = H; R10 = 4-ClC6H4; Z = OCH2C.tplbond.CH; X = O] (preparation given starting from 4-(2-aminoethyl)-2-methoxyphenol.HCl) with Lawesson's reagent afforded I [R1-R3 = H; R4 = Me; R5-R9 = H; R10 = 4-ClC6H4; Z = OCH2C.tplbond.CH; X = S] which inhibited fungal infestations by 80-100% at 200 ppm in tests against *Plasmopara viticola* on vines, *Phytophthora* on tomato plants, and *Phytophthora* on potato plants.

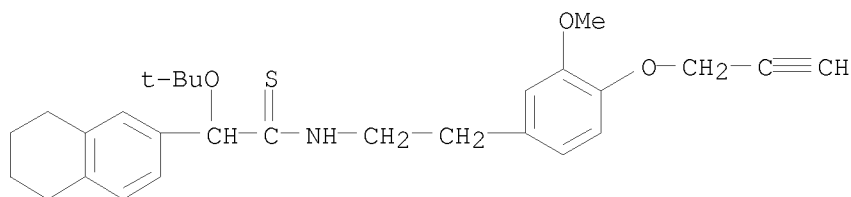
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	1106195-76-9		

RL: PRPH (Prophetic)

(Preparation of novel N-propargyloxyphenethyl thioacetamides as agrochemical fungicides)

RN 1106115-37-0 CAPLUS

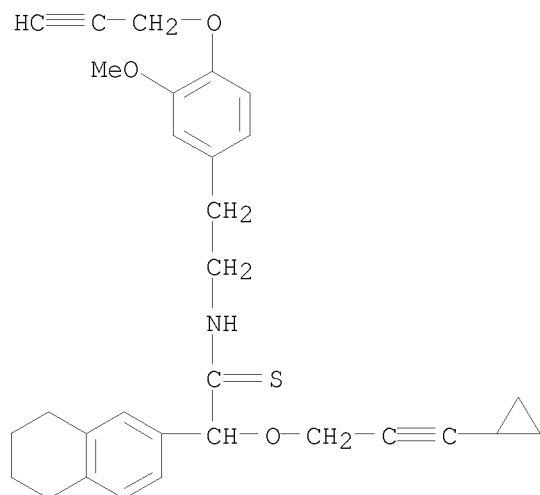
CN 2-Naphthaleneethanethioamide, α -(1,1-dimethylethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1106117-04-7 CAPLUS

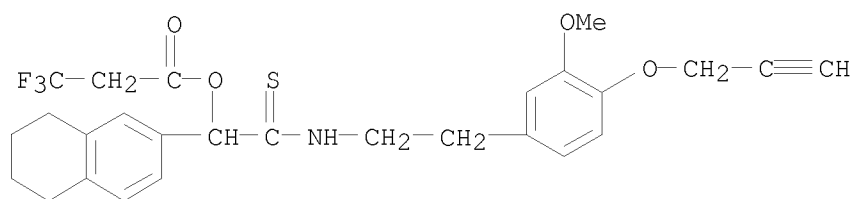
CN 2-Naphthaleneethanethioamide, α -[(3-cyclopropyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

10/513699



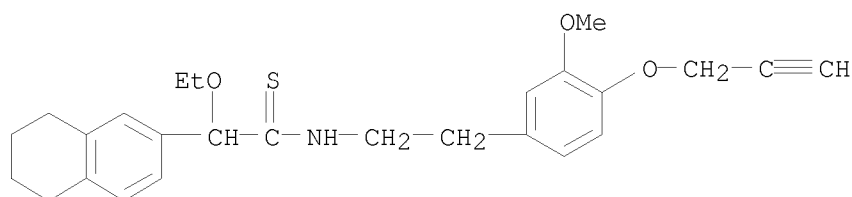
RN 1106118-72-2 CAPLUS

CN Propanoic acid, 3,3,3-trifluoro-, 2-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-1-(5,6,7,8-tetrahydro-2-naphthalenyl)-2-thioethoxy ester (CA INDEX NAME)



RN 1106119-83-8 CAPLUS

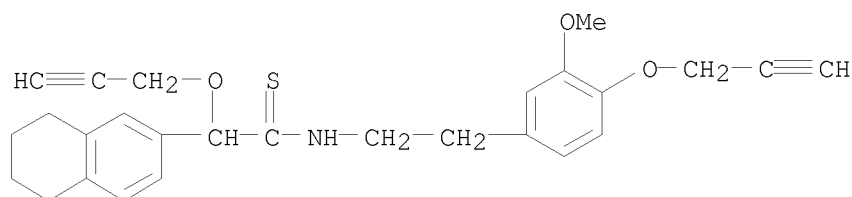
CN 2-Naphthaleneethanethioamide, α -ethoxy-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1106121-49-6 CAPLUS

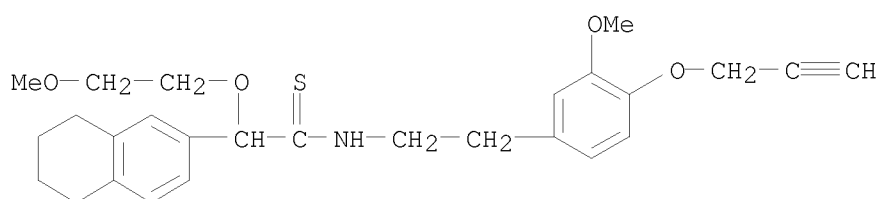
CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(2-propyn-1-yloxy)- (CA INDEX NAME)

10/513699



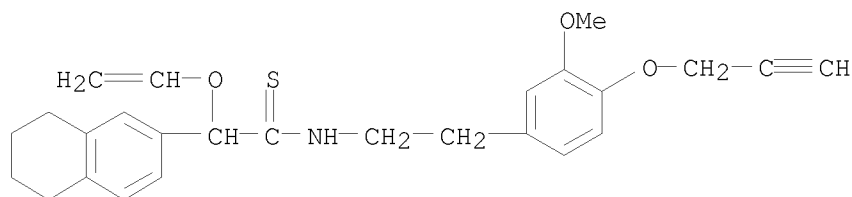
RN 1106123-14-1 CAPLUS

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RN 1106125-88-5 CAPLUS

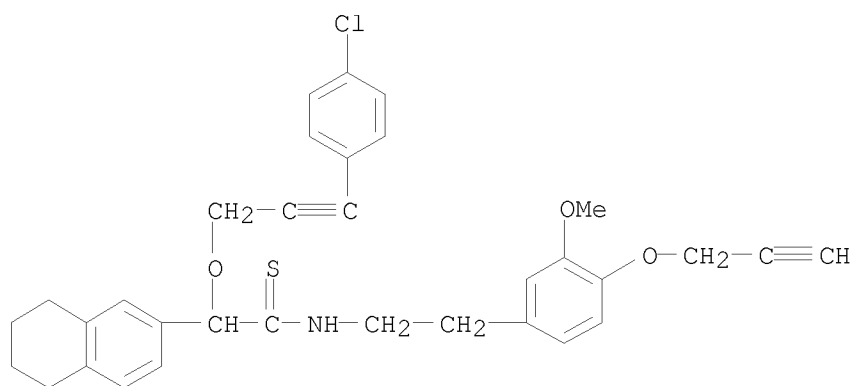
CN 2-Naphthaleneethanethioamide, α -(ethenyloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1106127-45-0 CAPLUS

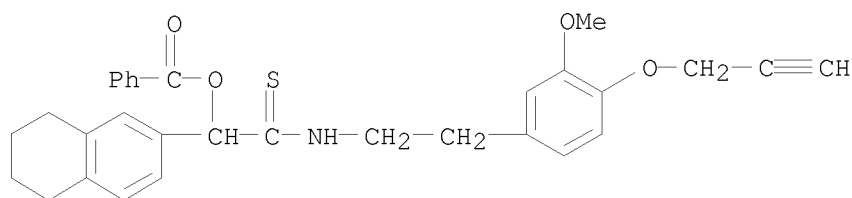
CN 2-Naphthaleneethanethioamide, α -[[3-(4-chlorophenyl)-2-propyn-1-yl]oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

10/513699



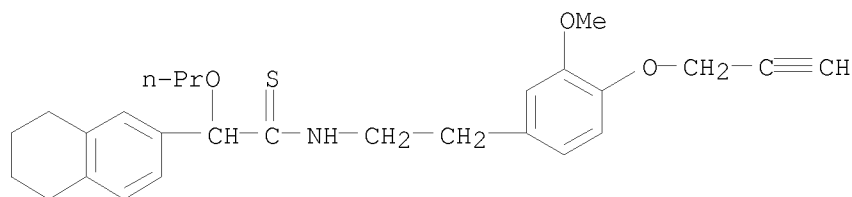
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CN 2-Naphthaleneethanethioamide, α -(benzoyloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1106130-24-8 CAPLUS

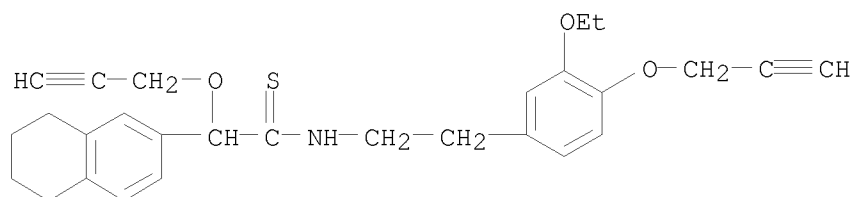
CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -propoxy- (CA INDEX NAME)



RN 1106131-88-7 CAPLUS

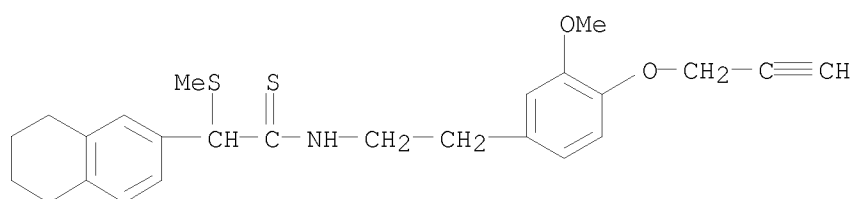
CN 2-Naphthaleneethanethioamide, N-[2-[3-ethoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)

10/513699



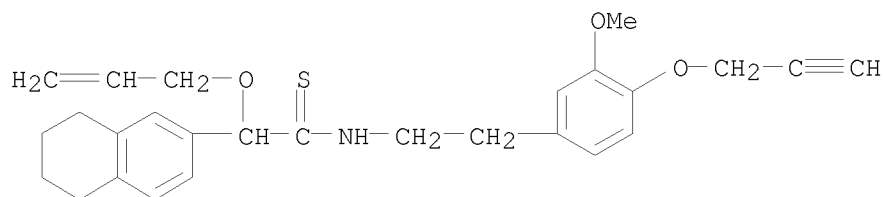
RN 1106133-55-4 CAPLUS

CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(methylthio)- (CA INDEX NAME)



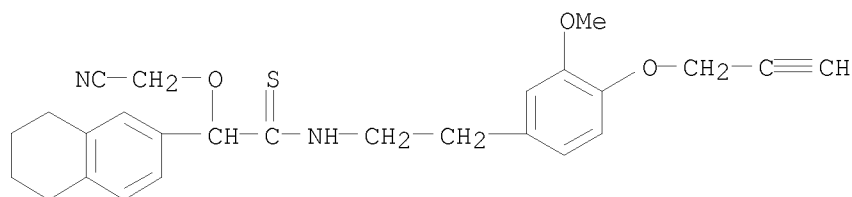
RN 1106140-59-3 CAPLUS

CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(2-propen-1-yloxy)- (CA INDEX NAME)



RN 1106142-26-0 CAPLUS

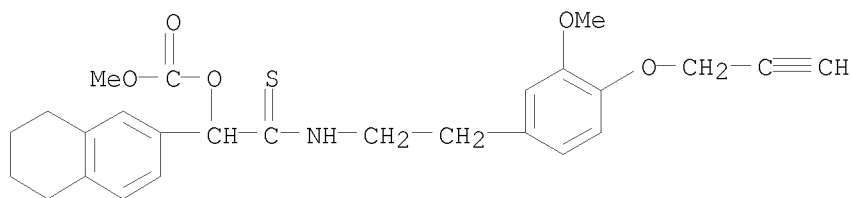
CN 2-Naphthaleneethanethioamide, α-(cyanomethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1106143-93-4 CAPLUS

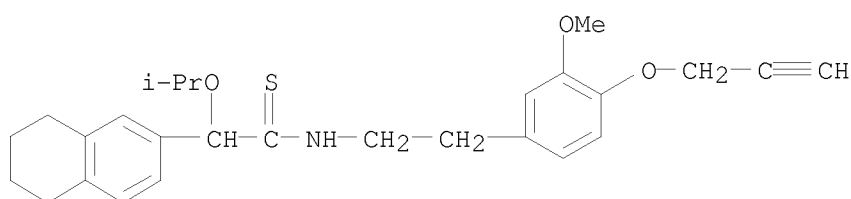
CN INDEX NAME NOT YET ASSIGNED

10/513699



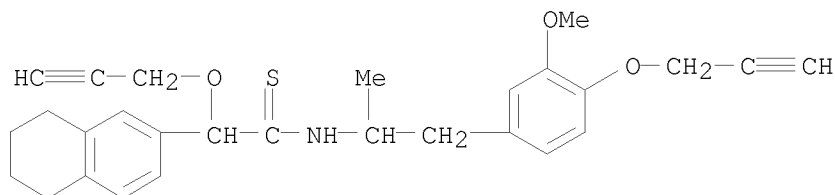
RN 1106145-05-4 CAPLUS

CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(1-methylethoxy)- (CA INDEX NAME)



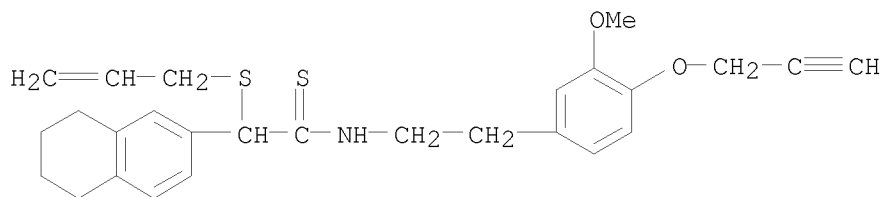
RN 1106146-72-8 CAPLUS

CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]-1-methylethyl]-α-(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1106148-39-3 CAPLUS

CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(2-propen-1-ylthio)- (CA INDEX NAME)

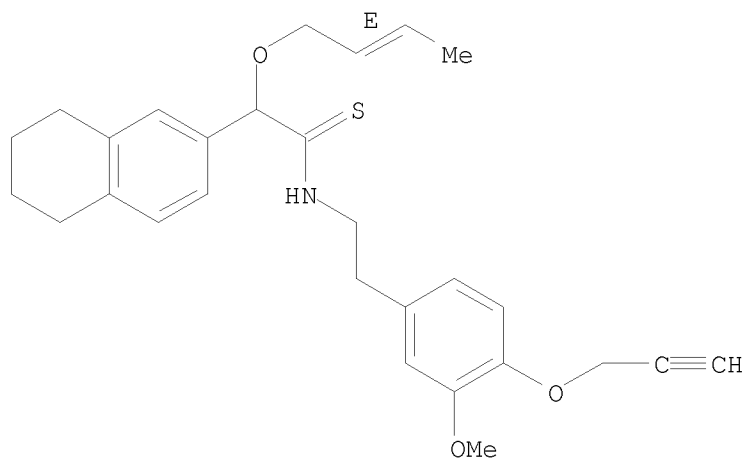


RN 1106151-18-1 CAPLUS

CN 2-Naphthaleneethanethioamide, α-[(2E)-2-buten-1-yloxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

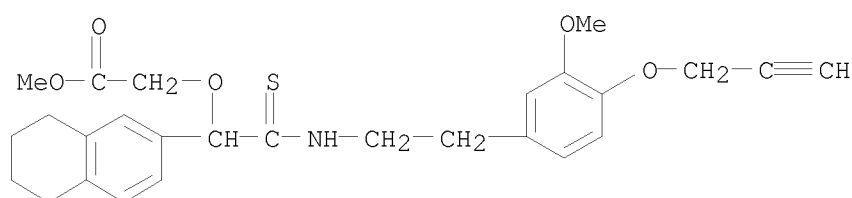
10/513699

Double bond geometry as shown.



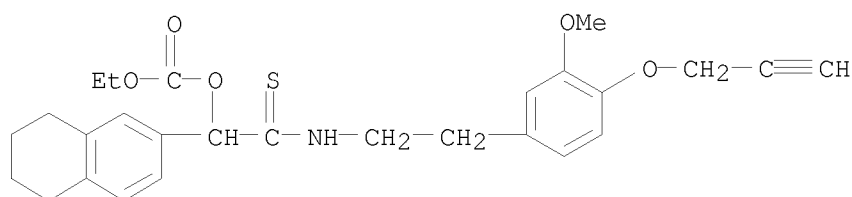
RN 1106152-85-5 CAPLUS

CN Acetic acid, 2-[2-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-1-(5,6,7,8-tetrahydro-2-naphthalenyl)-2-thioxoethoxy]-, methyl ester (CA INDEX NAME)



RN 1106154-52-2 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



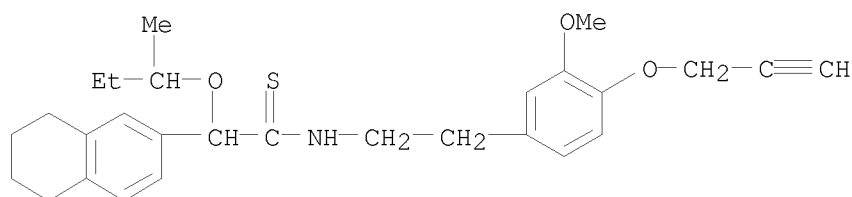
RN 1106155-64-9 CAPLUS

CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-alpha-(1-methylpropoxy)- (CA INDEX NAME)

<12/04/2007>

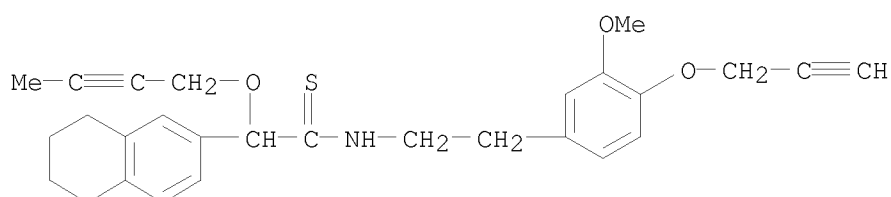
Erich Leese

10/513699



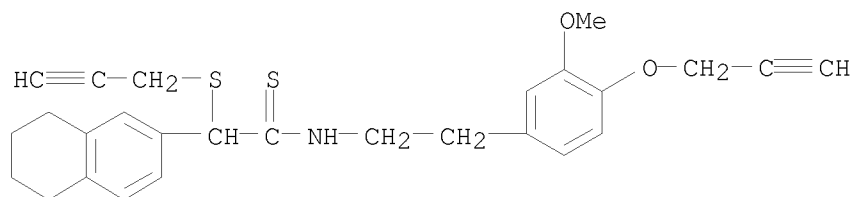
RN 1106157-31-6 CAPLUS

CN 2-Naphthaleneethanethioamide, α -(2-butyn-1-yloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



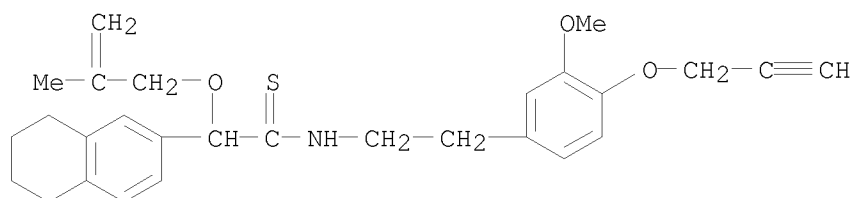
RN 1106158-98-8 CAPLUS

CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(2-propyn-1-ylthio)- (CA INDEX NAME)



RN 1106166-02-2 CAPLUS

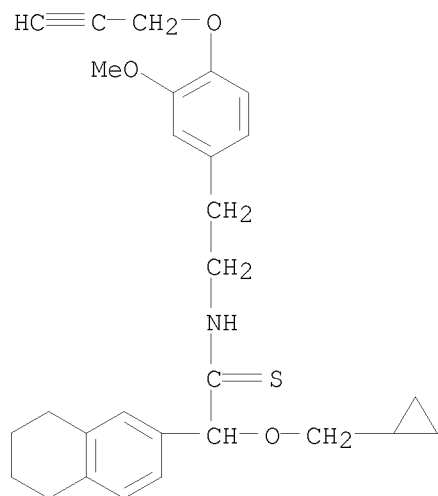
CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -[(2-methyl-2-propen-1-yl)oxy]- (CA INDEX NAME)



RN 1106167-69-4 CAPLUS

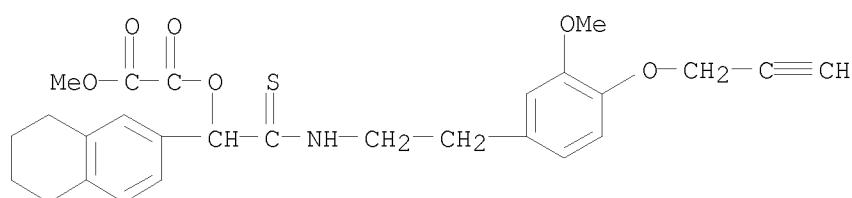
CN 2-Naphthaleneethanethioamide, α -(cyclopropylmethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

10/513699



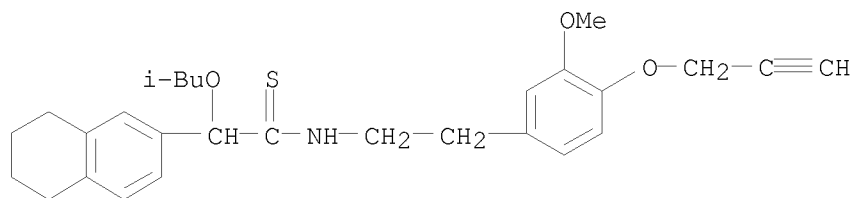
RN 1106169-36-1 CAPLUS

CN Ethanedioic acid, 1-[2-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-1-(5,6,7,8-tetrahydro-2-naphthalenyl)-2-thioxoethyl] 2-methyl ester (CA INDEX NAME)



RN 1106170-48-2 CAPLUS

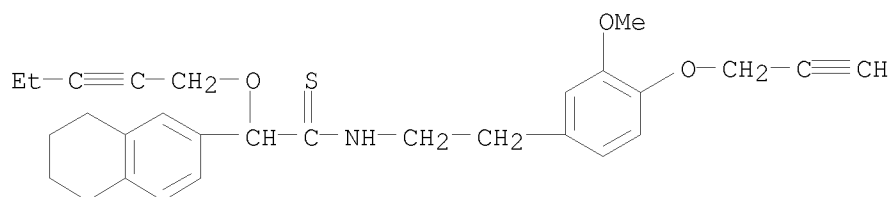
CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(2-methylpropoxy)- (CA INDEX NAME)



RN 1106172-14-8 CAPLUS

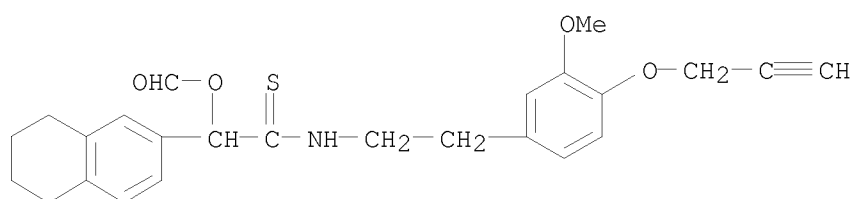
CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(2-pentyn-1-yloxy)- (CA INDEX NAME)

10/513699



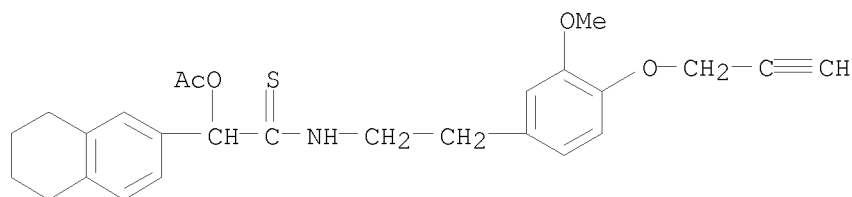
RN 1106173-78-7 CAPLUS

CN 2-Naphthaleneethanethioamide, α -(formyloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



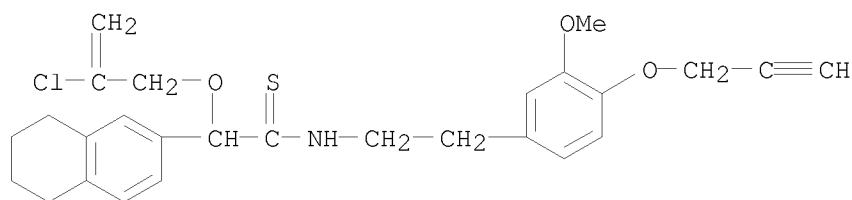
RN 1106175-32-9 CAPLUS

CN 2-Naphthaleneethanethioamide, α -(acetyloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1106176-99-1 CAPLUS

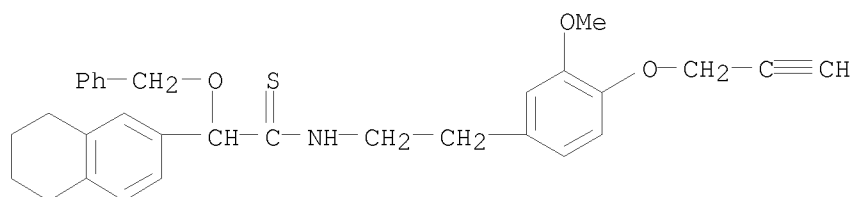
CN 2-Naphthaleneethanethioamide, α -[(2-chloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1106178-63-5 CAPLUS

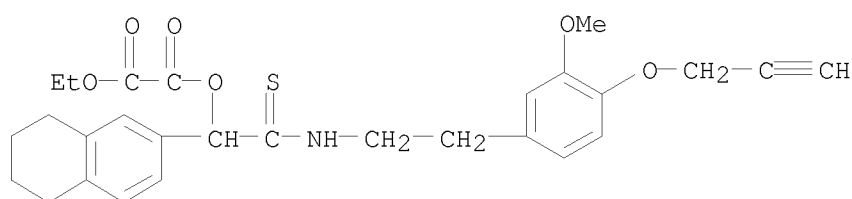
CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(phenylmethoxy)- (CA INDEX NAME)

10/513699



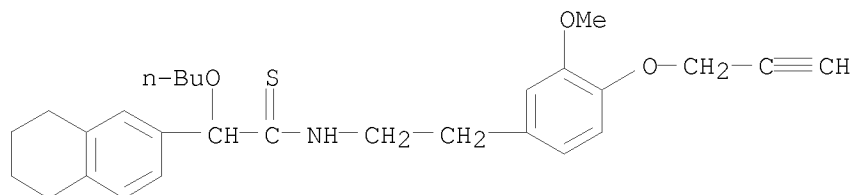
RN 1106180-29-3 CAPLUS

CN Ethanedioic acid, 1-ethyl 2-[2-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-1-(5,6,7,8-tetrahydro-2-naphthalenyl)-2-thioxoethyl] ester (CA INDEX NAME)



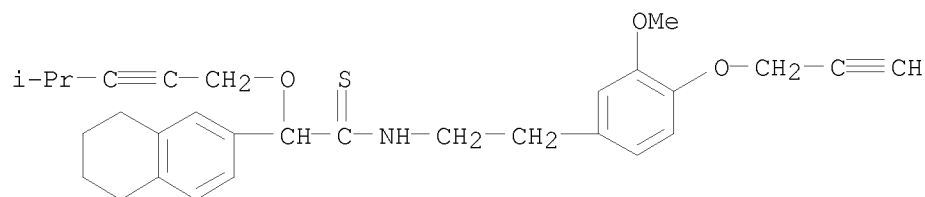
RN 1106182-20-0 CAPLUS

CN 2-Naphthaleneethanethioamide, α -butoxy-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1106183-87-2 CAPLUS

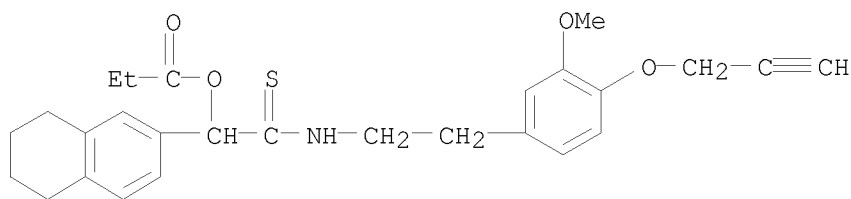
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RN 1106185-54-9 CAPLUS

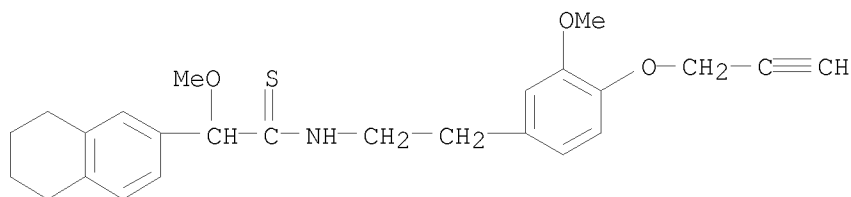
CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(1-oxopropoxy)- (CA INDEX NAME)

10/513699



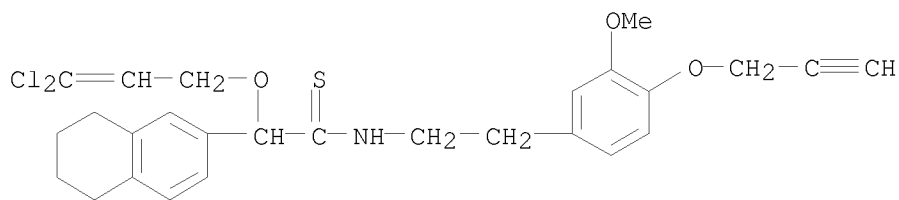
RN 1106192-49-7 CAPLUS

CN 2-Naphthaleneethanethioamide, 5,6,7,8-tetrahydro- α -methoxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



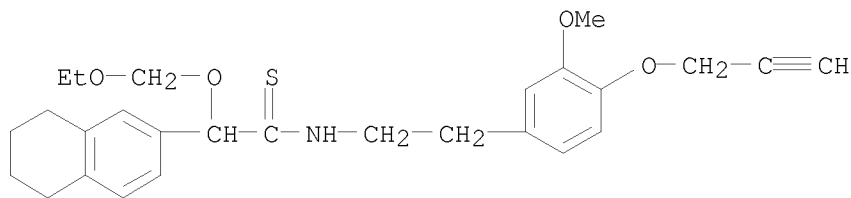
RN 1106194-10-8 CAPLUS

CN 2-Naphthaleneethanethioamide, α -[(3,3-dichloro-2-propen-1-yl)oxy]-
5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA
INDEX NAME)



RN 1106195-76-9 CAPLUS

CN 2-Naphthaleneethanethioamide, α -(ethoxymethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



REFERENCE COUNT:

2

THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L5 ANSWER 13 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2001:937747 CAPLUS

DOCUMENT NUMBER: 136:410929

TITLE: Preparation of ureidoalkylpiperidines as modulators of chemokine CCR3 receptor activity.

INVENTOR(S): Ko, Soo S.; Delucca, George V.; Duncia, John V.; Santella, Joseph B.; Wacker, Dean A.; Yao, Wenqing

PATENT ASSIGNEE(S): Dupont Pharmaceuticals Company, USA; Bristol-Myers Squibb Pharmaceutical Co.

SOURCE: PCT Int. Appl., 446 pp.

CODEN: PIXXD2

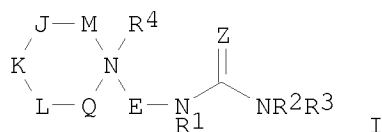
DOCUMENT TYPE: Patent

LANGUAGE: English

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2001098269 A2		20011227	WO 2001-XI19745	20010620
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW				
RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG, TR				
PRIORITY APPLN. INFO.:			US 2000-213051P	20000621
			US 2000-598821	20000621

GI



AB [Title compds. I; M = CH₂, CHR₅, CHR₁₃, CR₁₃R₁₃, CR₅R₁₃; Q = CH₂, CHR₅, CHR₁₃, CR₁₃R₁₃, CR₅R₁₃; J, L = CH₂, CHR₅, CHR₆, CR₆R₆, CR₅R₆; Z = O, S; M = CH₂, CHR₅, CHR₁₃, CR₁₃R₁₃, CR₅R₁₃; K = CHR₅, CR₅R₆; Z = O, S; E = (CHR₇)(CHR₉)v(CR₁₁R₁₂); R₁, R₂ = H, alkyl, alkenyl, alkynyl, (substituted) alkylcycloalkyl; R₂R₃ = atoms to form a (substituted) 5-7 membered ring; R₃, R₅ = (substituted) (alkyl)cycloalkyl, (alkyl)heterocyclyl; R₄ = null, O, alkyl, alkenyl, alkynyl, etc.; R₄ with R₇, R₉, or R₁₁ = atoms to form a 5-7 membered ring; R₇, R₉ = H; R₄R₇, R₄R₉ = (substituted) spirocyclyl; R₁₃ = alkyl, alkenyl, alkynyl, cycloalkyl, etc.; R₁₁R₁₂ = pyrrolidinyl, tetrahydrofuryl, piperidinyl, tetrahydropyranyl; v = 1, 2], were prepared as modulators of chemokine activity (no data). Thus, 4-benzyl-1-(3-aminopropyl)piperidine (preparation given) in THF was treated with 3-cyanophenyl isocyanate to give N-(3-cyanophenyl)-N'-[3-[4-(phenylmethyl)-1-piperidinyl]propyl]urea. [This abstract record is one of 15 records for this document necessitated by the large number of index entries required to fully index the document and publication system constraints.]

IT 1084141-39-8

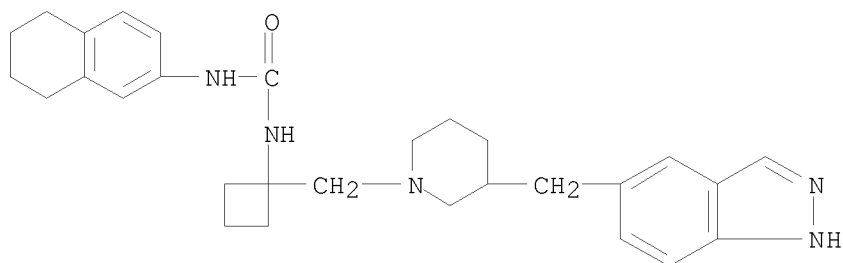
RL: PRPH (Prophetic)

10/513699

(Preparation of ureidoalkylpiperidines as modulators of chemokine CCR3
receptor activity.)

RN 1084141-39-8 CAPLUS

CN Urea, N-[1-[[3-(1H-indazol-5-ylmethyl)-1-piperidinyl]methyl]cyclobutyl]-N'-
(5,6,7,8-tetrahydro-2-naphthalenyl)- (CA INDEX NAME)



REFERENCE COUNT:

5

THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L5 ANSWER 14 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 2001:851100 CAPLUS

DOCUMENT NUMBER: 135:371520

TITLE: Preparation of novel phenyl propargyl ethers as agrochemical fungicides

INVENTOR(S): Lamberth, Clemens; Zeller, Martin; Kunz, Walter; Cederbaum, Fredrik

PATENT ASSIGNEE(S): Syngenta Participations A.-G., Switz.

SOURCE: PCT Int. Appl., 84 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

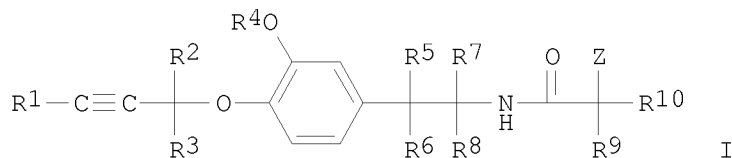
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AU 2001060301	A	20011126	AU 2001-60301	20010515 <--
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HU 2003001965	A3	20090128		
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AT 271031	T	20040715	AT 2001-933967	20010515 <--
ES 2223848	T3	20050301	ES 2001-933967	20010515
RU 2259353	C2	20050827	RU 2002-133216	20010515
CN 1221526	C	20051005	CN 2001-809580	20010515
AP 1601	A	20060430	AP 2002-2649	20010515
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IN 2002CN01841	A	20050211	IN 2002-CN1841	20021111
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ZA 2002009266	A	20031020	ZA 2002-9266	20021114 <--
US 6683211	B1	20040127	US 2002-276476	20021115 <--
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			WO 2001-EO5530	W 20010515
			WO 2001-EP5530	W 20010515

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 135:371520

GI



AB The title compds. [I; R1 = H, alkyl, cycloalkyl, (un)substituted aryl; R2, R3 = H, alkyl; R4 = alkyl, alkenyl, alkynyl; R5- R8 = H, alkyl; R9 = H, (un)substituted alkyl, alkenyl or alkynyl; R10 = (un)substituted (hetero)aryl; Z = halo, (un)substituted aryloxy, alkoxy, etc.] which possess useful plant protecting properties and may advantageously be employed in agricultural practice for controlling or preventing the infestation of plants by phytopathogenic microorganisms, especially fungi (biol.

data given), were prepared E.g., a multi-step synthesis of I [R1-R3 = H; R4 = Me; R5-R8 = H; R9 = H; R10 = 4-ClC6H4; Z = OMe] was given.

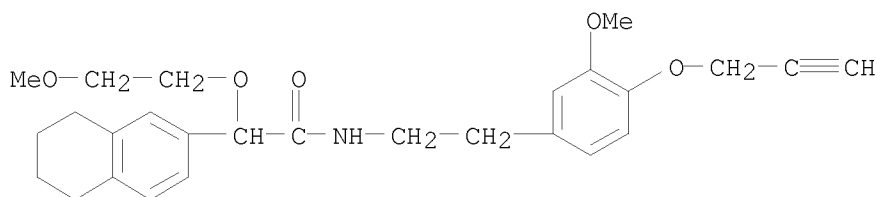
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RL: PRPH (Prophetic)

(Preparation of novel phenyl propargyl ethers as agrochemical fungicides)

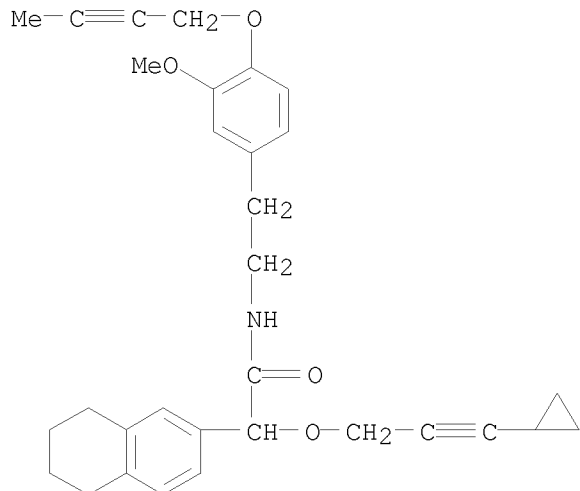
RN 1055179-98-0 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -(2-methoxyethoxy)-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

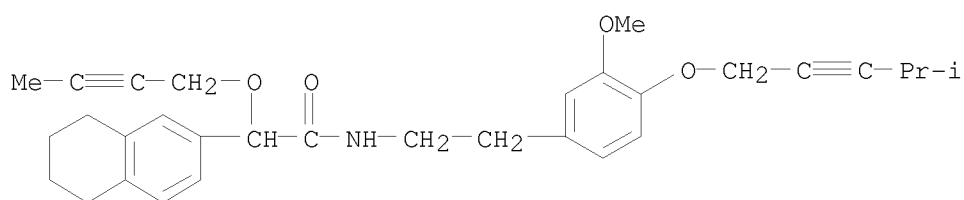
RN 1055181-48-0 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -(3-cyclopropyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro- (CA INDEX NAME)

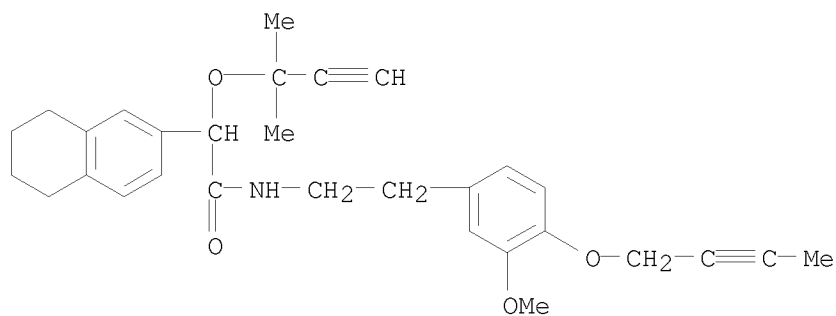
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RN 1055182-12-1 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

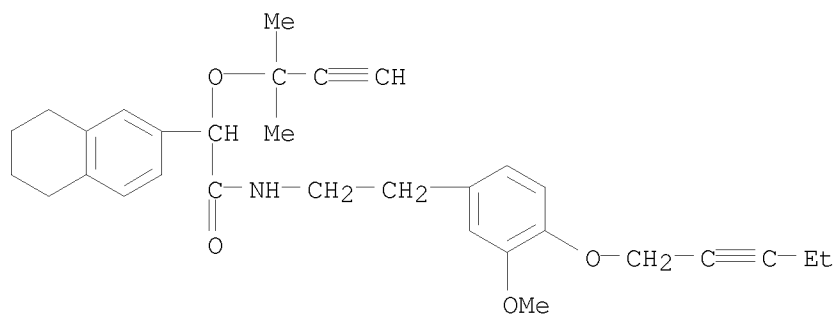


RN 1055182-57-4 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-
α-[(1,1-dimethyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro- (CA INDEX
NAME)

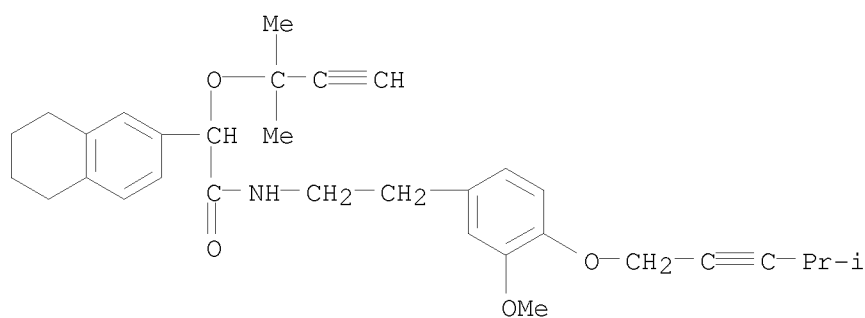


RN 1055182-58-5 CAPLUS
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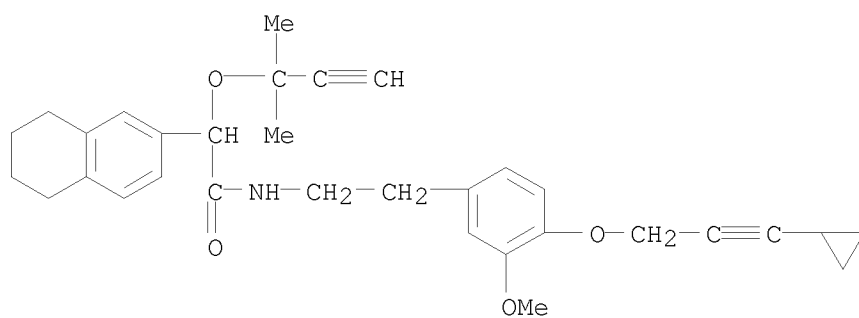
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RN 1055182-59-6 CAPLUS
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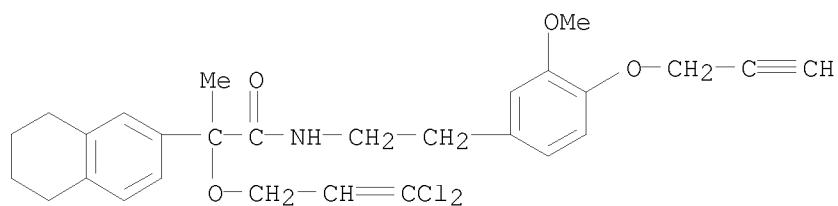


RN 1055182-60-9 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]- α -[(1,1-dimethyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro- (CA INDEX NAME)



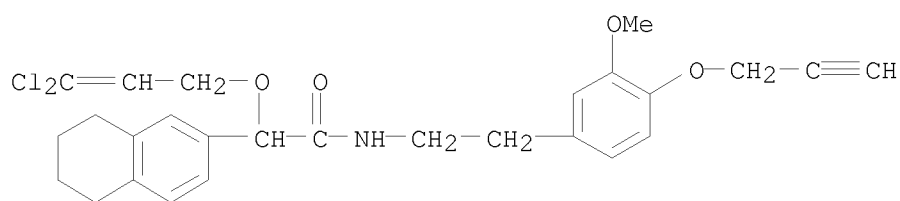
RN 1055183-36-2 CAPLUS
CN 2-Naphthaleneacetamide, α -[(3,3-dichloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)

10/513699



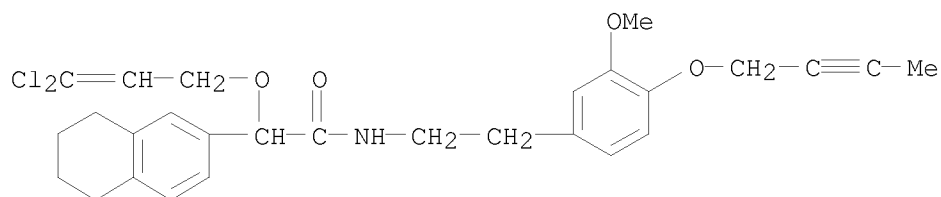
RN 1055183-37-3 CAPLUS

CN 2-Naphthaleneacetamide, α -[(3,3-dichloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



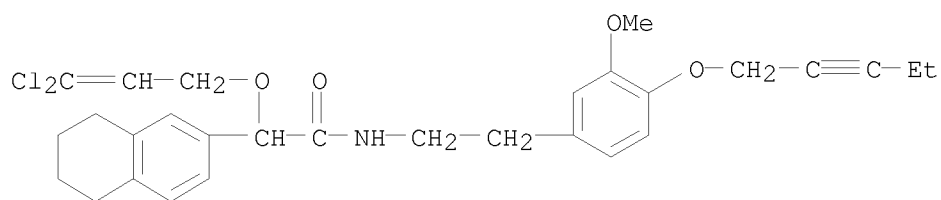
RN 1055183-38-4 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -[(3,3-dichloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 1055183-39-5 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



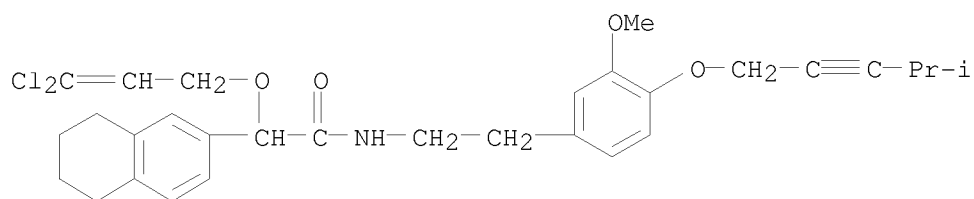
RN 1055183-40-8 CAPLUS

CN INDEX NAME NOT YET ASSIGNED

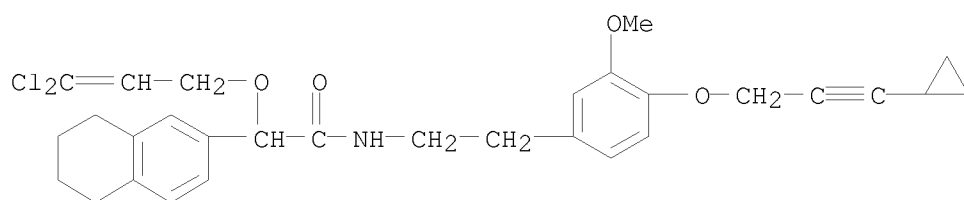
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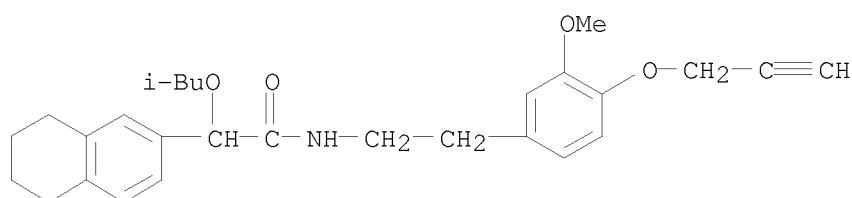
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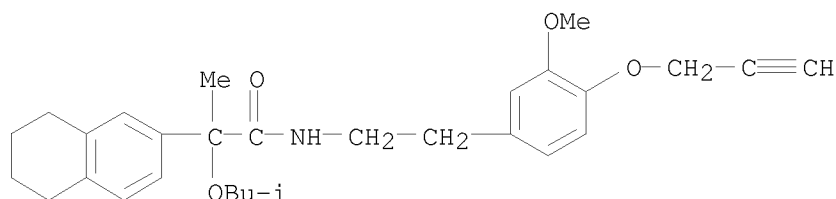
RN 1055183-41-9 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055185-17-5 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(2-methylpropoxy)- (CA INDEX NAME)

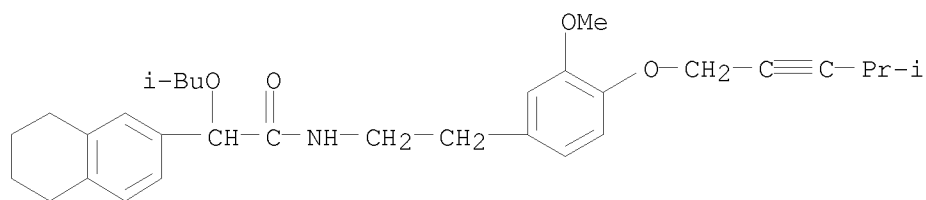


RN 1055185-18-6 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- α -(2-methylpropoxy)- (CA INDEX NAME)



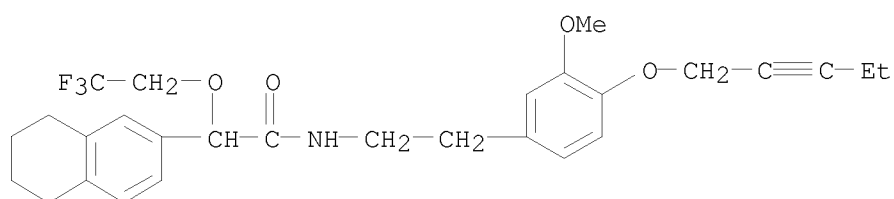
RN 1055185-19-7 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

10/513699



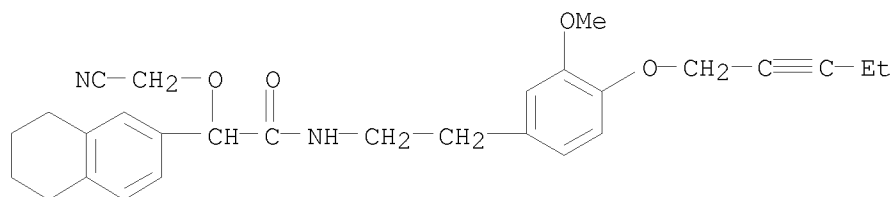
RN 1055186-14-5 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]-α-(2,2,2-trifluoroethoxy)- (CA INDEX NAME)



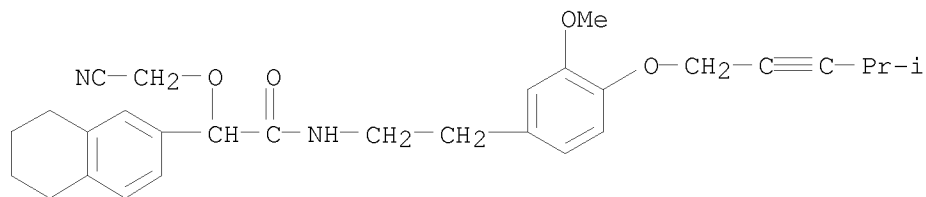
RN 1055186-58-7 CAPLUS

CN 2-Naphthaleneacetamide, α-(cyanomethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055186-59-8 CAPLUS

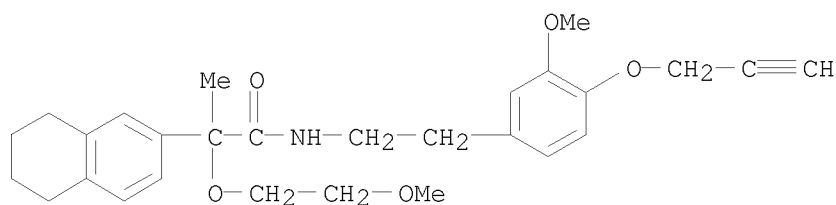
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RN 1055189-32-6 CAPLUS

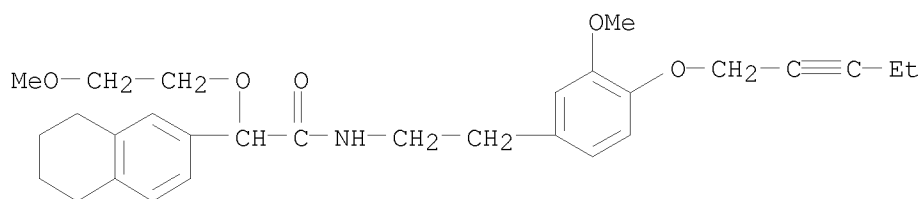
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-α-(2-methoxyethoxy)-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl- (CA INDEX NAME)

10/513699



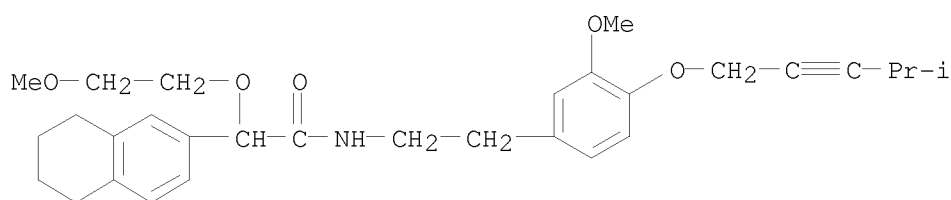
RN 1055189-33-7 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -(2-methoxyethoxy)-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



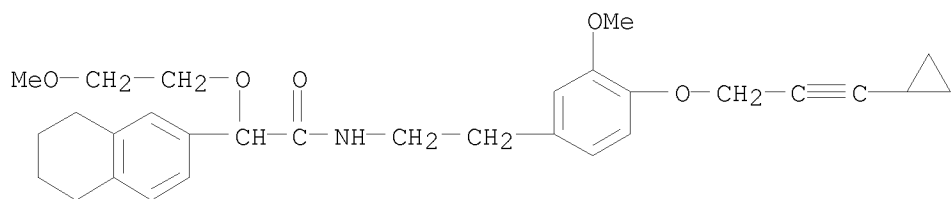
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CN INDEX NAME NOT YET ASSIGNED



RN 1055189-35-9 CAPLUS

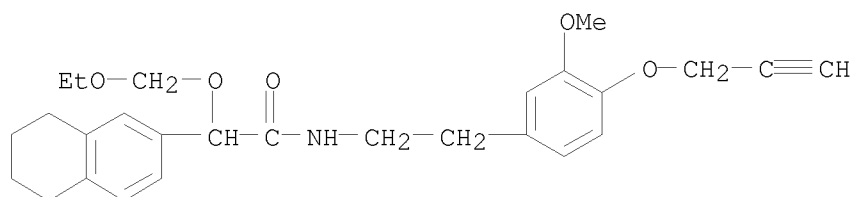
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-methoxyethoxy)- (CA INDEX NAME)



RN 1055199-07-9 CAPLUS

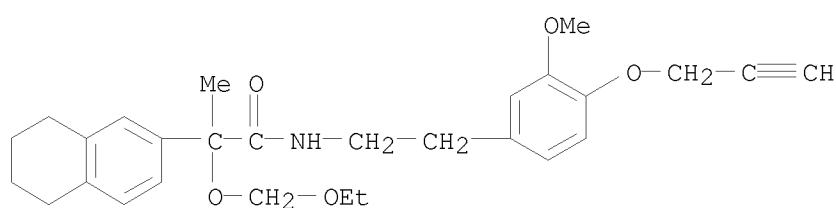
CN 2-Naphthaleneacetamide, α -(ethoxymethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

10/513699



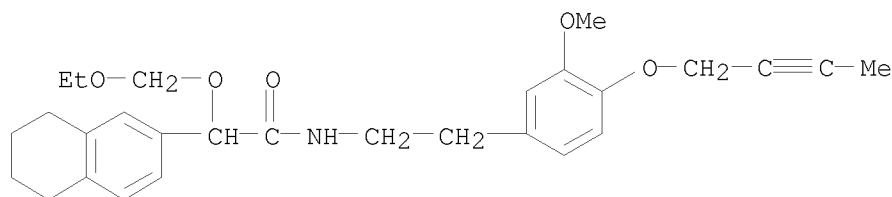
RN 1055199-08-0 CAPLUS

CN 2-Naphthaleneacetamide, α -(ethoxymethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



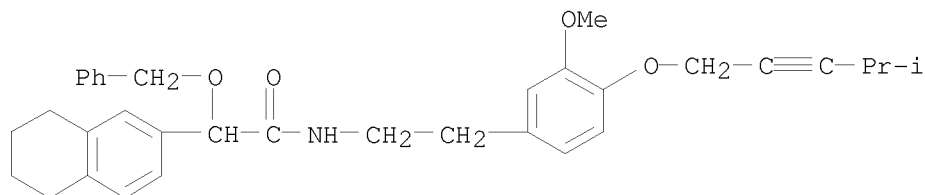
RN 1055199-09-1 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -(ethoxymethoxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)



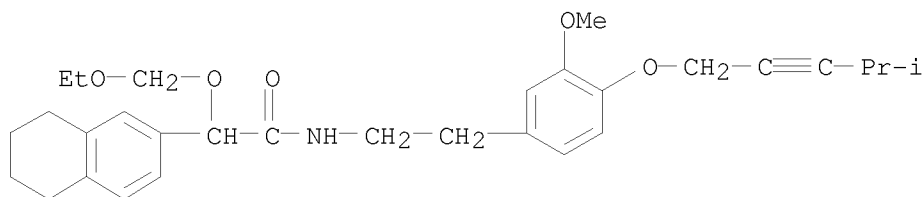
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CN INDEX NAME NOT YET ASSIGNED

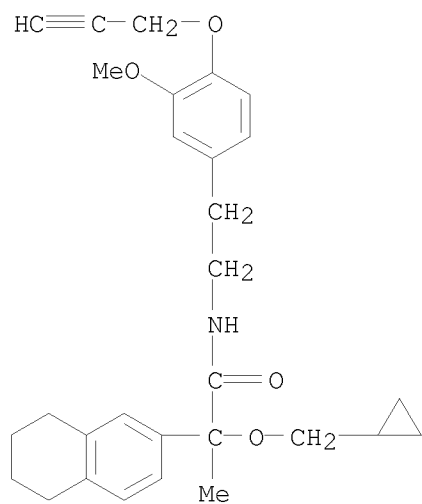


RN 1055203-63-8 CAPLUS

CN 2-Naphthaleneacetamide, α -(ethoxymethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

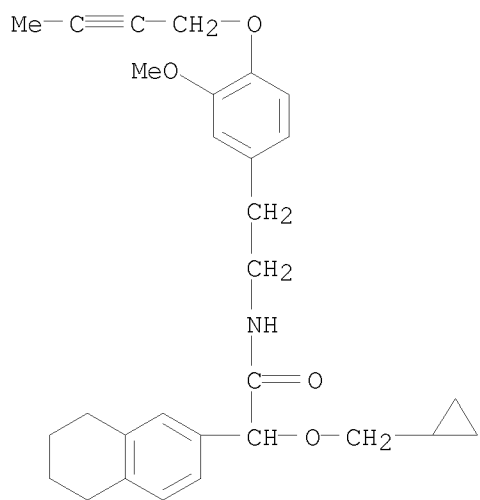
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10/513699



RN 1055206-26-2 CAPLUS

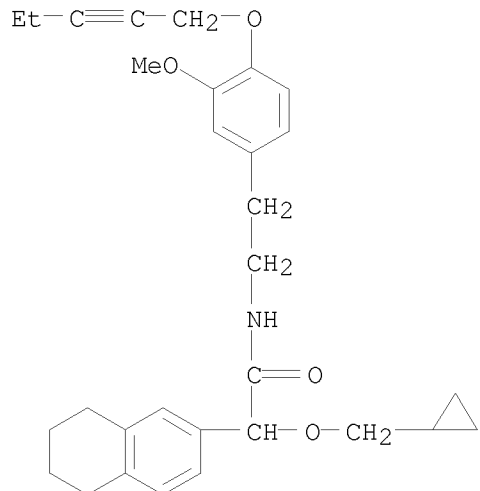
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-
α-(cyclopropylmethoxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)



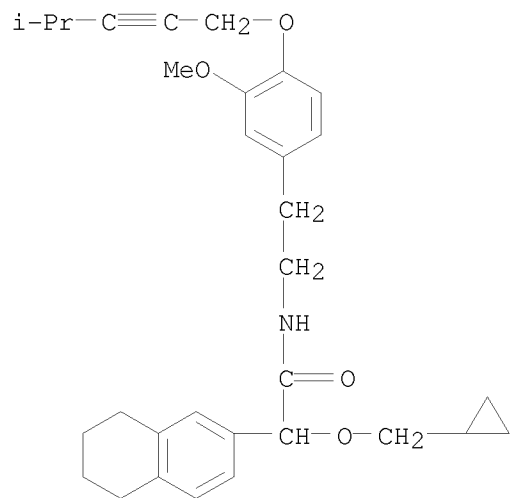
RN 1055206-27-3 CAPLUS

CN 2-Naphthaleneacetamide, α-(cyclopropylmethoxy)-5,6,7,8-tetrahydro-N-
[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

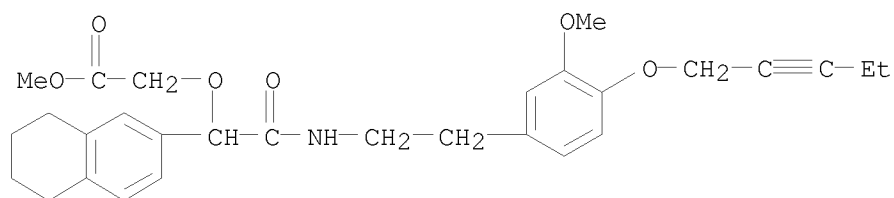
10/513699



RN 1055206-28-4 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055207-50-5 CAPLUS
CN Acetic acid, 2-[2-[[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]amino]-2-oxo-1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethoxy]-, methyl ester (CA INDEX NAME)



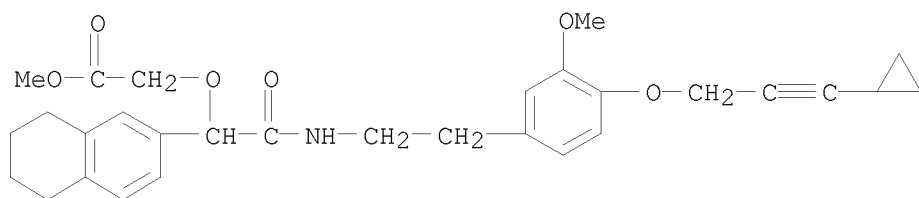
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10/513699

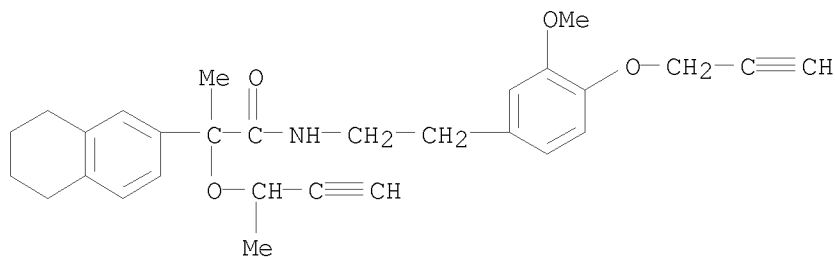
RN 1055207-51-6 CAPLUS

CN Acetic acid, 2-[2-[[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]amino]-2-oxo-1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethoxy]-, methyl ester (CA INDEX NAME)



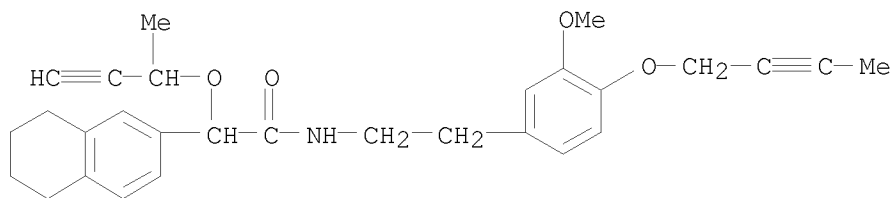
RN 1055208-05-3 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- α -[(1-methyl-2-propyn-1-yl)oxy]- (CA INDEX NAME)



RN 1055208-06-4 CAPLUS

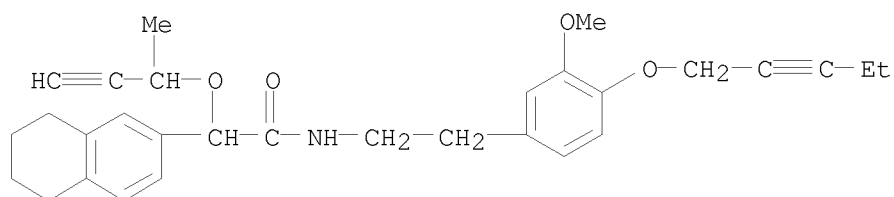
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -[(1-methyl-2-propyn-1-yl)oxy]- (CA INDEX NAME)



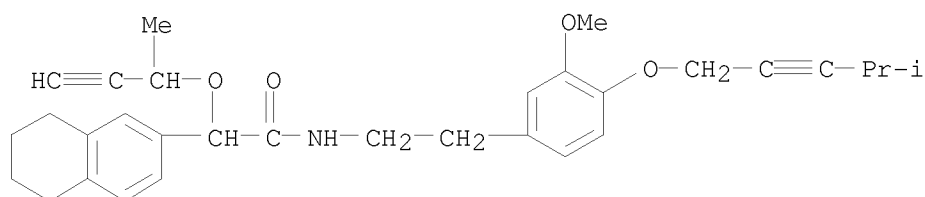
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CN INDEX NAME NOT YET ASSIGNED

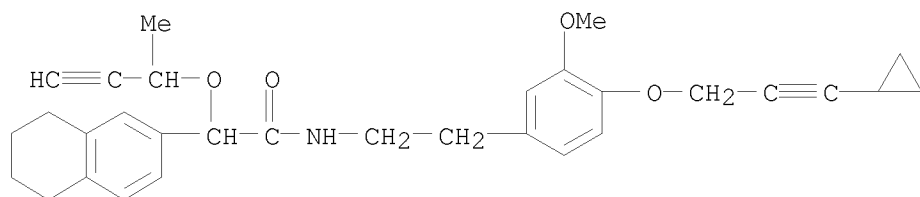
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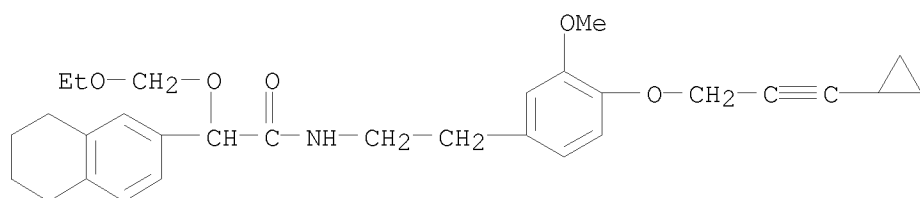
RN 1055208-08-6 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055208-09-7 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-[(1-methyl-2-propyn-1-yl)oxy]- (CA INDEX NAME)

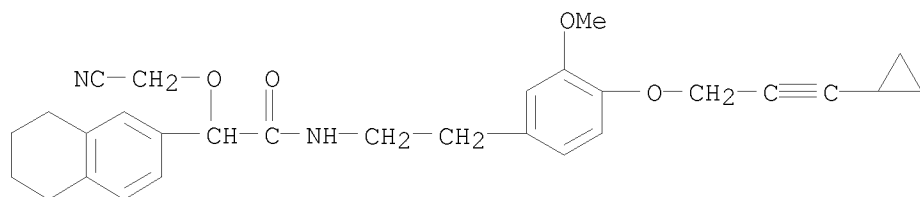


RN 1055208-72-4 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-α-(ethoxymethoxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)

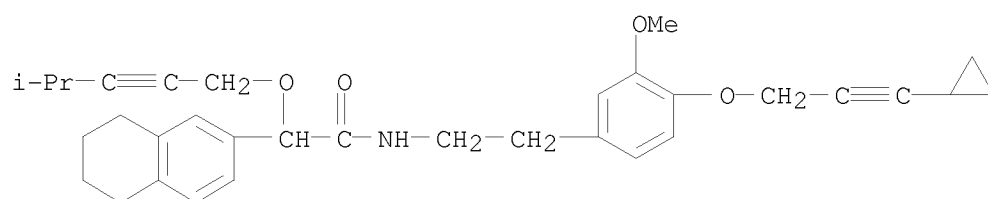


RN 1055209-45-4 CAPLUS
CN 2-Naphthaleneacetamide, α-(cyanomethoxy)-N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)

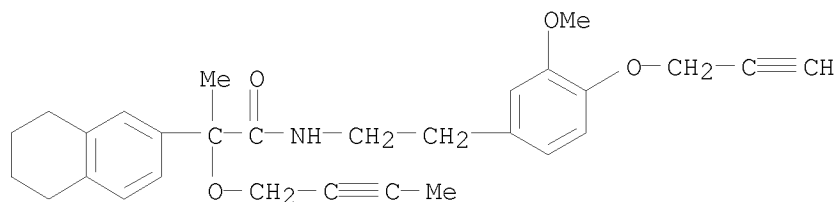
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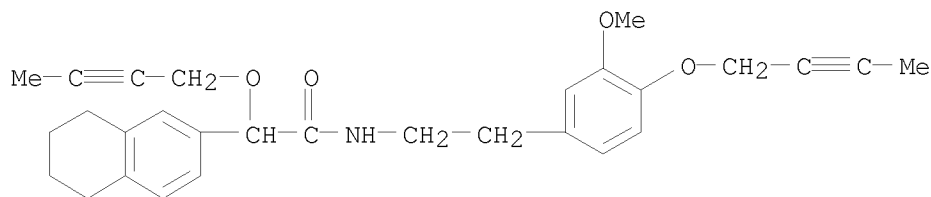
RN 1055210-27-9 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055214-37-3 CAPLUS
CN 2-Naphthaleneacetamide, α -(2-butyn-1-yloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)

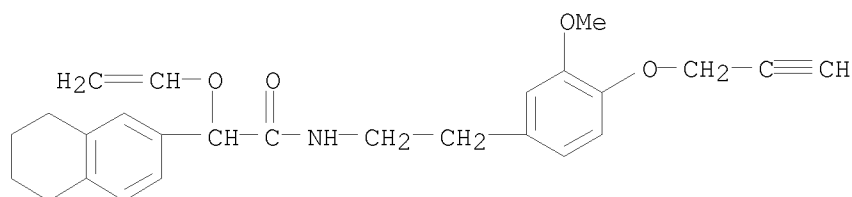


RN 1055214-38-4 CAPLUS
CN 2-Naphthaleneacetamide, α -(2-butyn-1-yloxy)-N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)



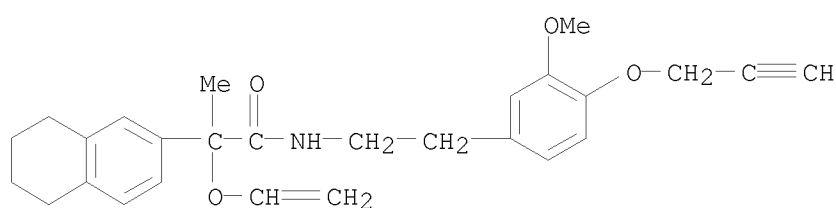
RN 1055215-71-8 CAPLUS
CN 2-Naphthaleneacetamide, α -(ethenyloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

10/513699



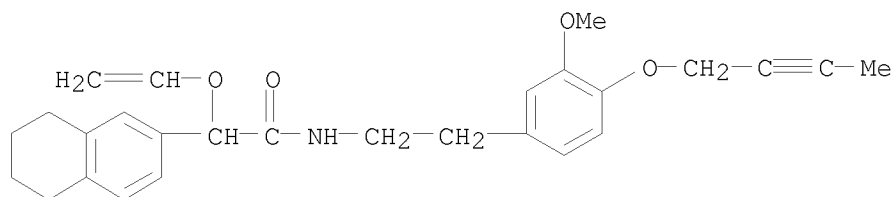
RN 1055215-72-9 CAPLUS

CN 2-Naphthaleneacetamide, α -(ethenyloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



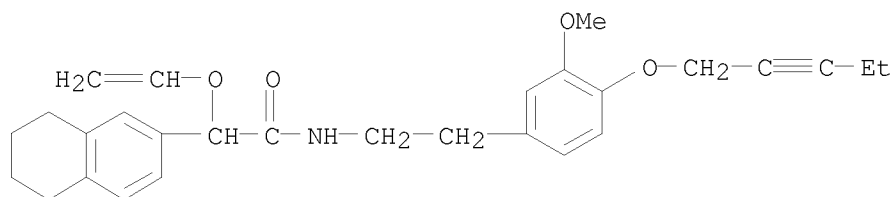
RN 1055215-73-0 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -(ethenyloxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 1055215-74-1 CAPLUS

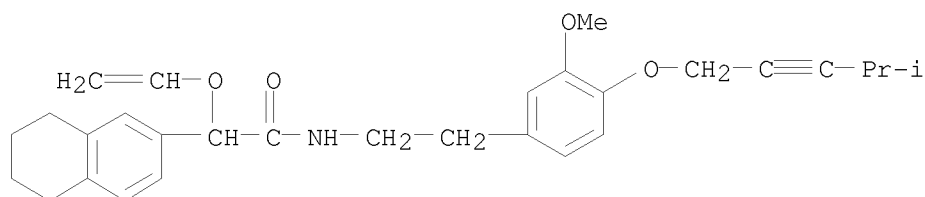
CN 2-Naphthaleneacetamide, α -(ethenyloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055215-75-2 CAPLUS

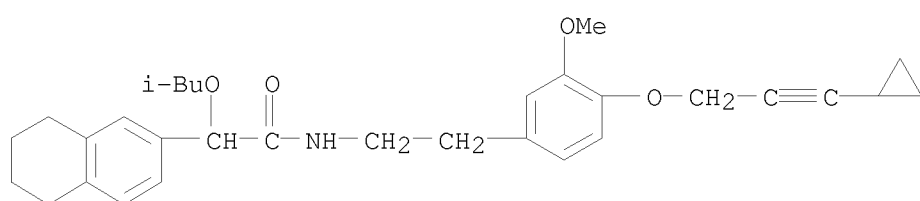
CN INDEX NAME NOT YET ASSIGNED

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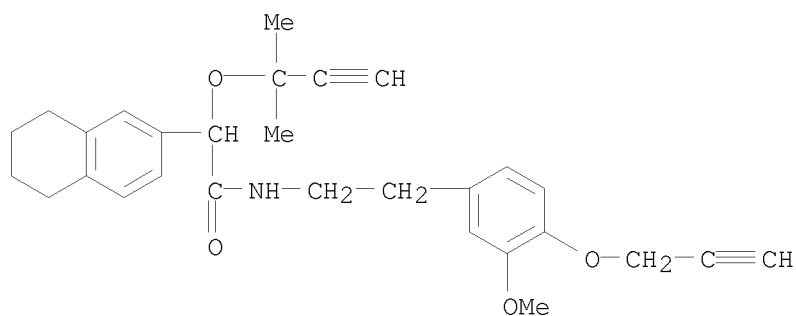
RN 1055216-06-2 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(2-methylpropoxy)- (CA INDEX NAME)



RN 1055216-39-1 CAPLUS

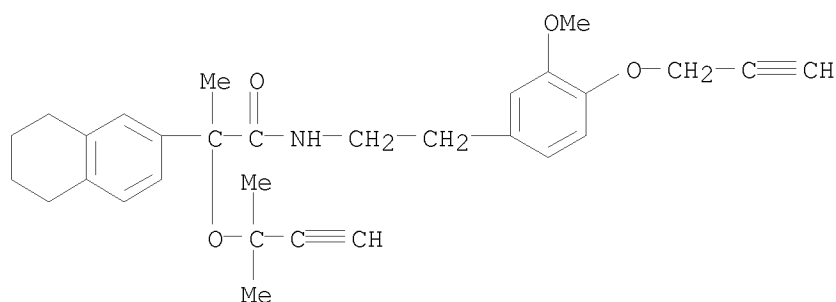
CN 2-Naphthaleneacetamide, α-[(1,1-dimethyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055216-40-4 CAPLUS

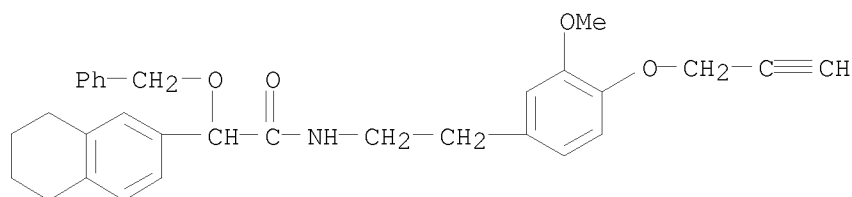
CN 2-Naphthaleneacetamide, α-[(1,1-dimethyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl- (CA INDEX NAME)

10/513699



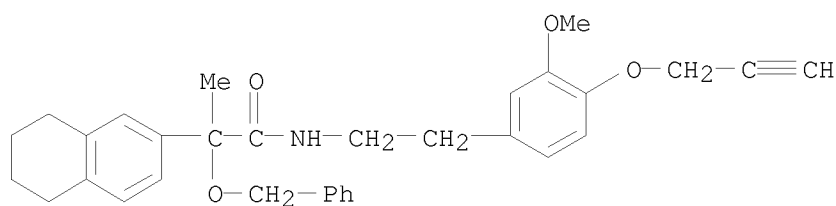
RN 1055216-75-5 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(phenylmethoxy)- (CA INDEX NAME)



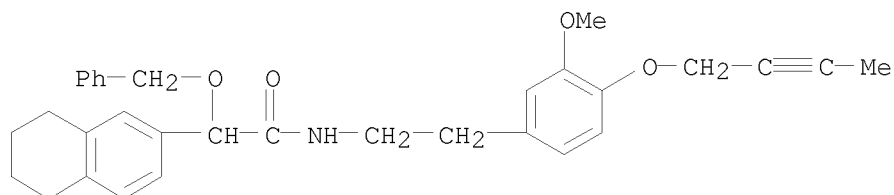
RN 1055216-76-6 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-(phenylmethoxy)- (CA INDEX NAME)



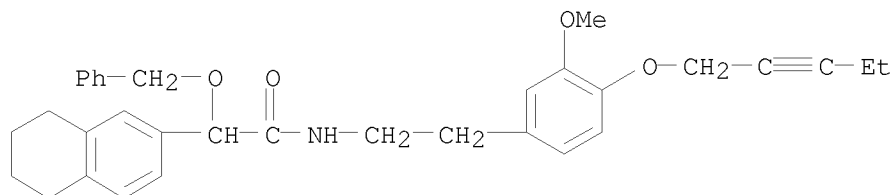
RN 1055216-77-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(phenylmethoxy)- (CA INDEX NAME)

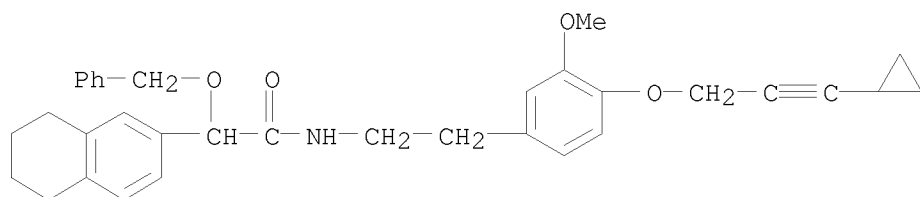


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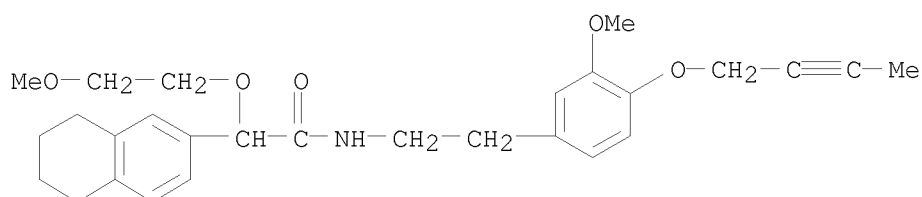
RN 1055216-78-8 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- α -(phenylmethoxy)- (CA INDEX NAME)



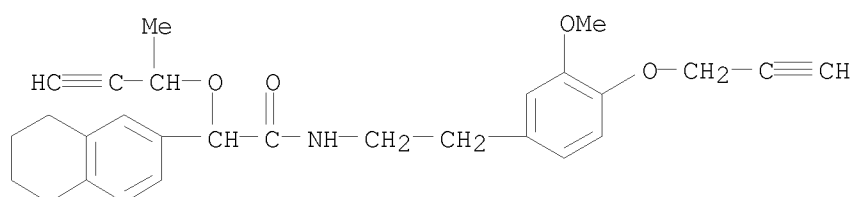
RN 1055216-79-9 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(phenylmethoxy)- (CA INDEX NAME)



RN 1055216-94-8 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-methoxyethoxy)- (CA INDEX NAME)



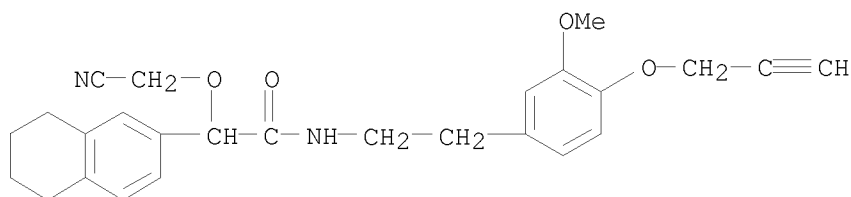
RN 1055217-05-4 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(1-methyl-2-propyn-1-yloxy)- (CA INDEX NAME)



10/513699

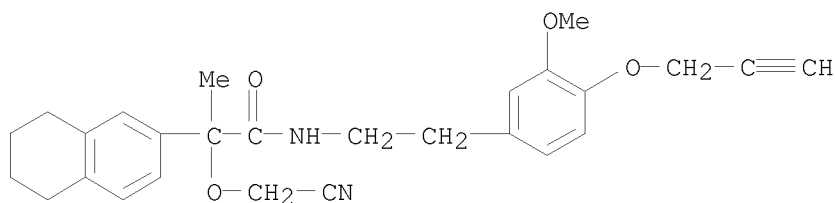
RN 1055217-88-3 CAPLUS

CN 2-Naphthaleneacetamide, α -(cyanomethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



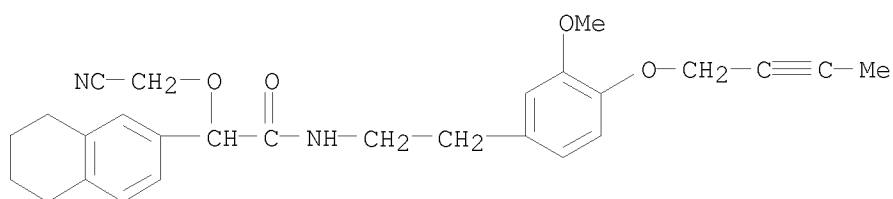
RN 1055217-89-4 CAPLUS

CN 2-Naphthaleneacetamide, α -(cyanomethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



RN 1055217-90-7 CAPLUS

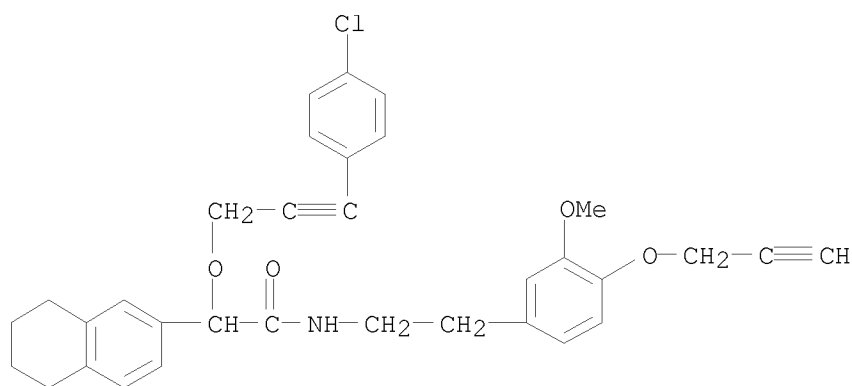
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -(cyanomethoxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 1055218-09-1 CAPLUS

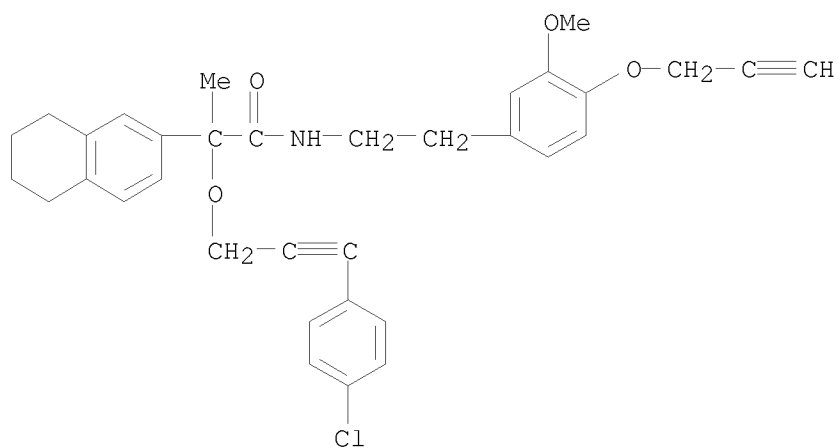
CN 2-Naphthaleneacetamide, α -[[3-(4-chlorophenyl)-2-propyn-1-yl]oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

10/513699



RN 1055218-10-4 CAPLUS

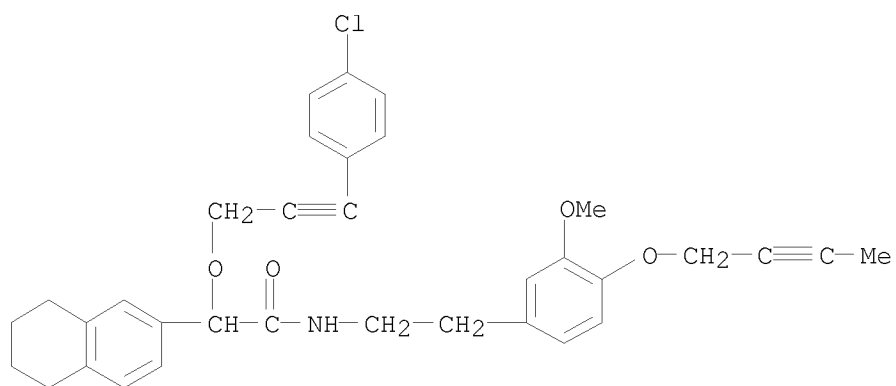
CN 2-Naphthaleneacetamide, α -[[3-(4-chlorophenyl)-2-propyn-1-yl]oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



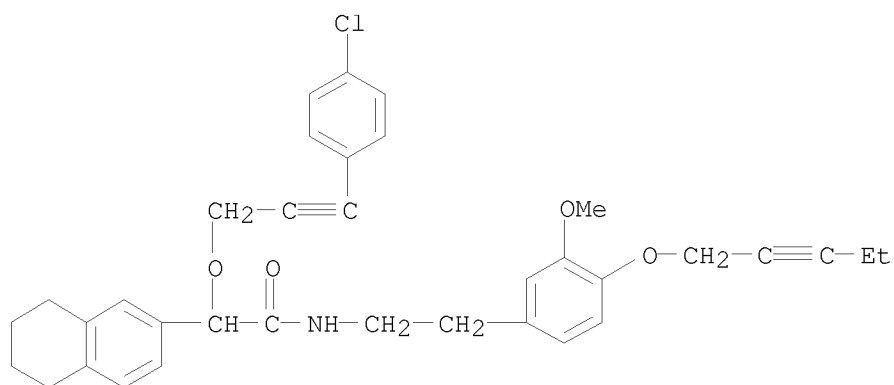
RN 1055218-11-5 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -[[3-(4-chlorophenyl)-2-propyn-1-yl]oxy]-5,6,7,8-tetrahydro- (CA INDEX NAME)

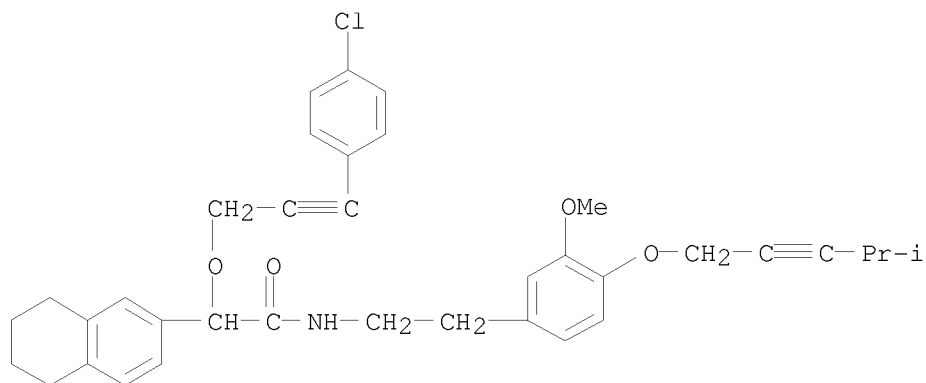
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RN 1055218-12-6 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

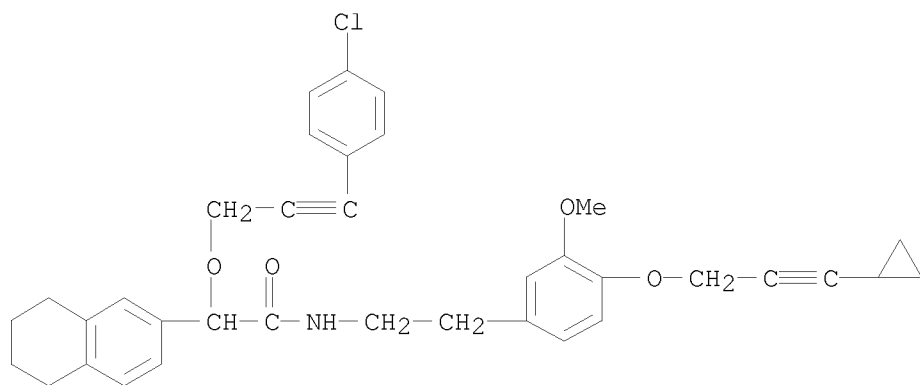


RN 1055218-13-7 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

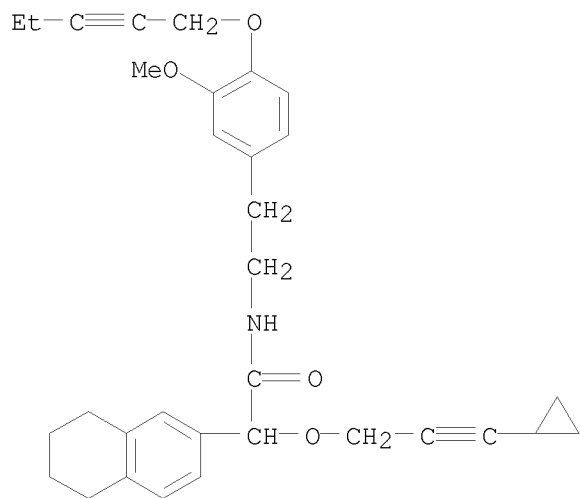


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RN 1055218-14-8 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

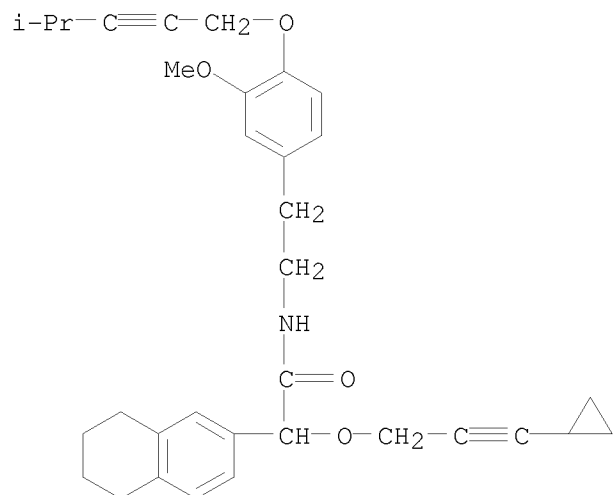


RN 1055219-85-6 CAPLUS
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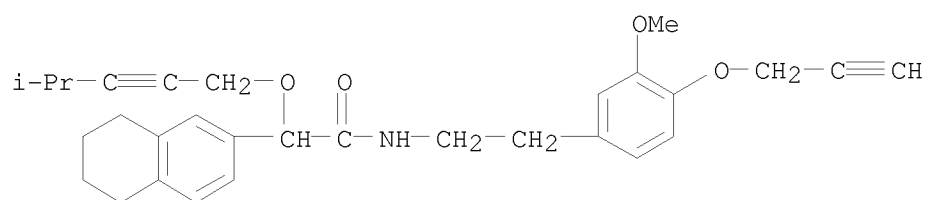


RN 1055219-86-7 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

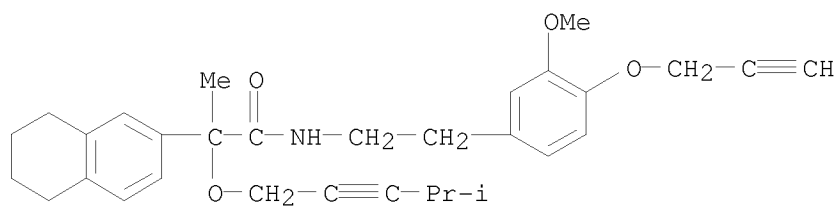
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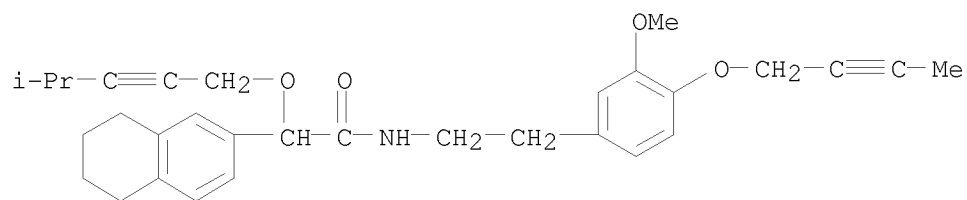


RN 1055220-20-6 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

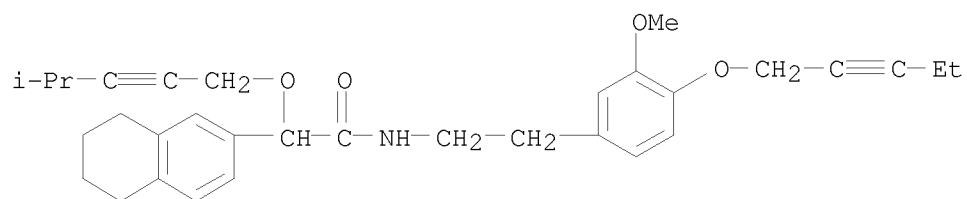


RN 1055220-21-7 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

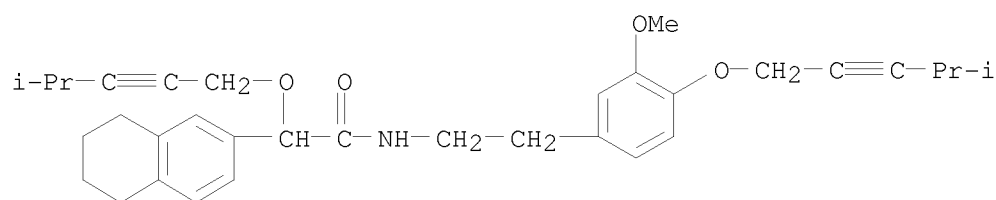
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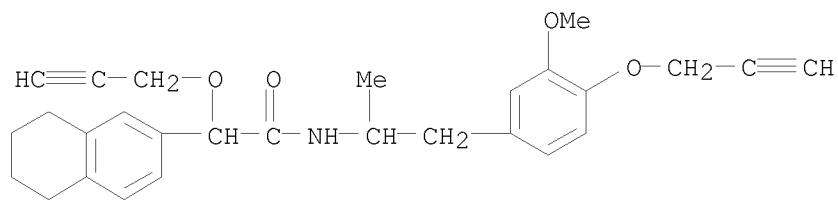
RN 1055220-22-8 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055220-23-9 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

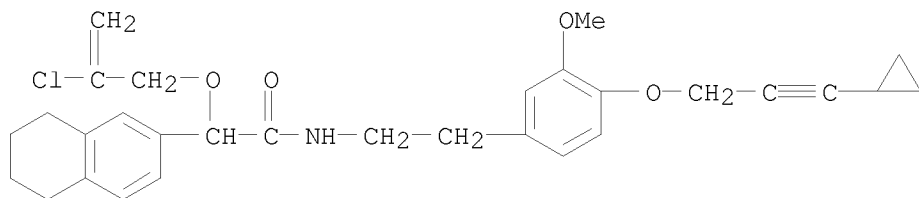


RN 1055221-97-0 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]-1-methylethyl]-alpha-(2-propyn-1-yloxy)- (CA INDEX NAME)



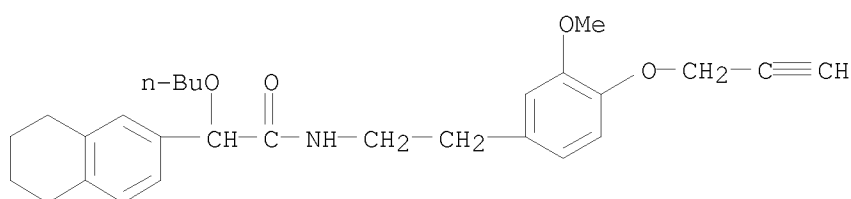
RN 1055222-32-6 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

10/513699



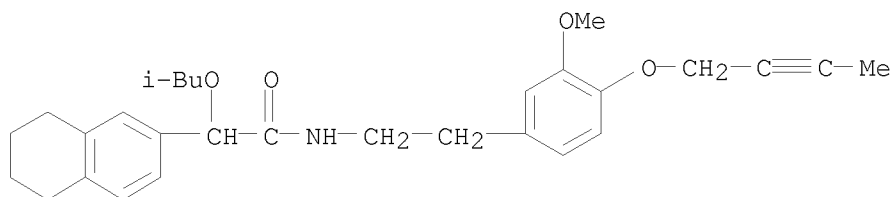
RN 1055223-53-4 CAPLUS

CN 2-Naphthaleneacetamide, α -butoxy-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



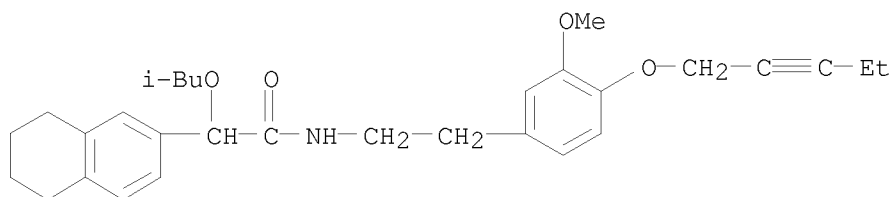
RN 1055224-57-1 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-methylpropoxy)- (CA INDEX NAME)



RN 1055224-58-2 CAPLUS

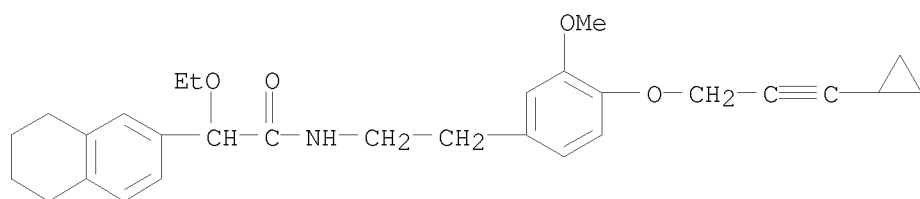
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- α -(2-methylpropoxy)- (CA INDEX NAME)



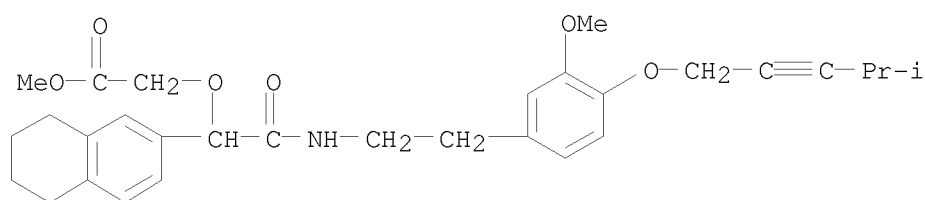
RN 1055225-51-8 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]- α -ethoxy-5,6,7,8-tetrahydro- (CA INDEX NAME)

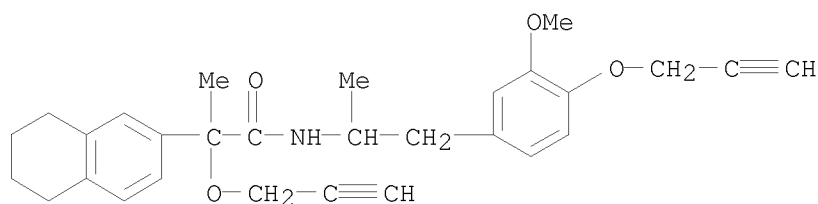
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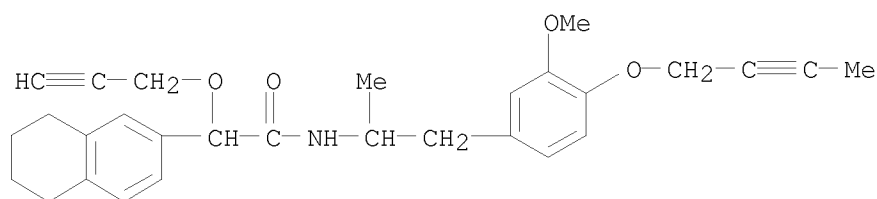
RN 1055226-57-7 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055228-55-1 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]-1-methylethyl]-alpha-methyl-alpha-(2-propyn-1-yloxy)- (CA INDEX NAME)

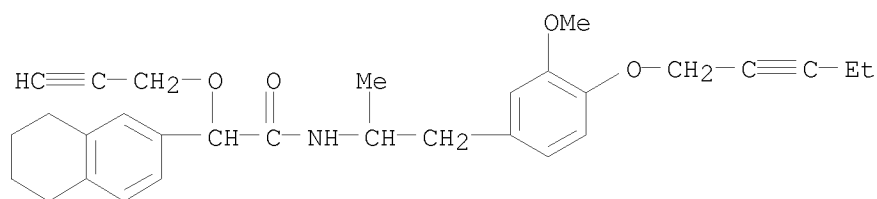


RN 1055228-56-2 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]-1-methylethyl]-5,6,7,8-tetrahydro-alpha-(2-propyn-1-yloxy)- (CA INDEX NAME)

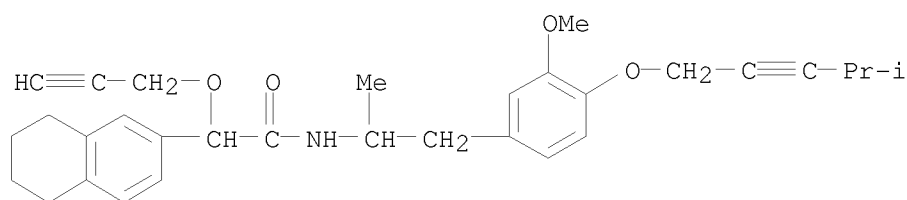


RN 1055228-57-3 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]-1-methylethyl]-alpha-(2-propyn-1-yloxy)- (CA INDEX NAME)

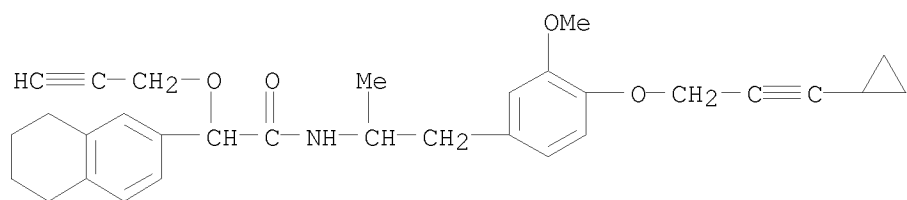
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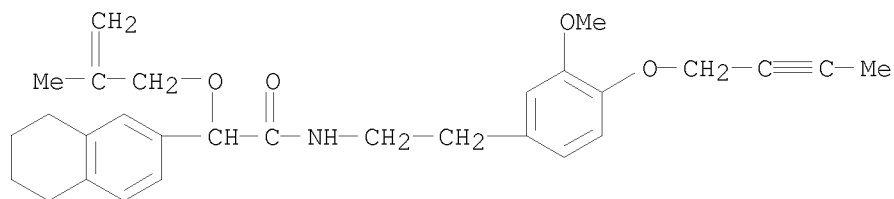
RN 1055228-58-4 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055228-59-5 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]-1-methylethyl]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)

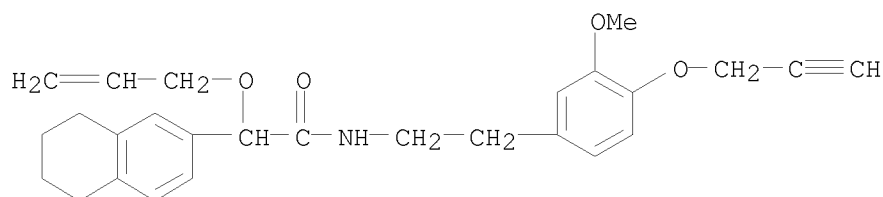


RN 1055229-27-0 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -[(2-methyl-2-propen-1-yl)oxy]- (CA INDEX NAME)

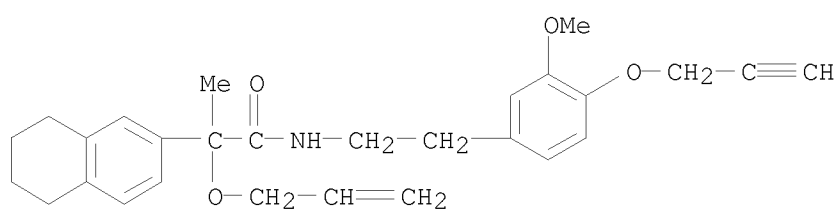


RN 1055230-40-4 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(2-propen-1-yloxy)- (CA INDEX NAME)

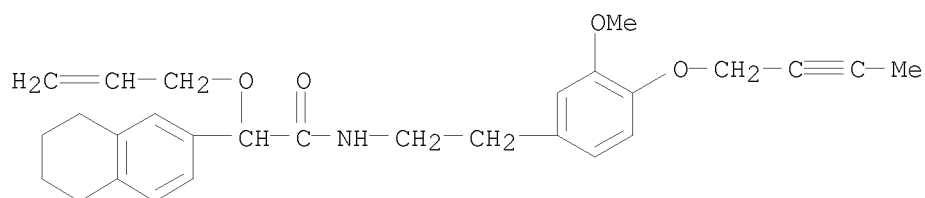
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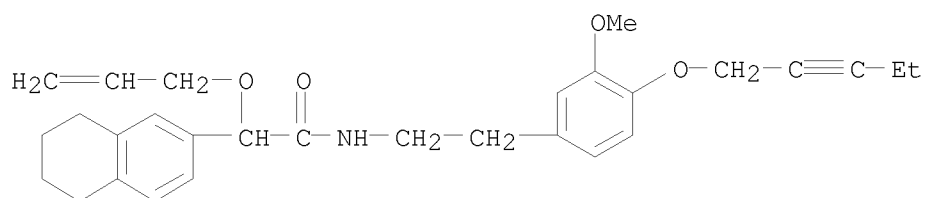
RN 1055230-41-5 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- α -(2-propen-1-yloxy)- (CA INDEX NAME)



RN 1055230-42-6 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-propen-1-yloxy)- (CA INDEX NAME)

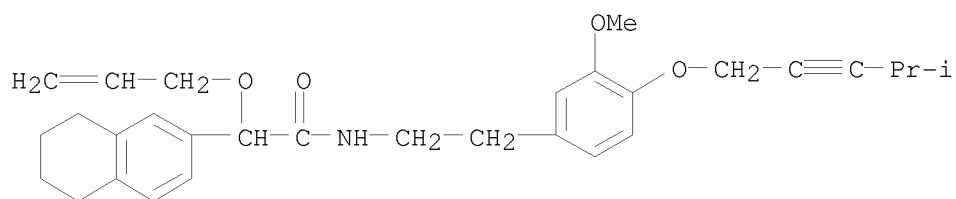


RN 1055230-43-7 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- α -(2-propen-1-yloxy)- (CA INDEX NAME)



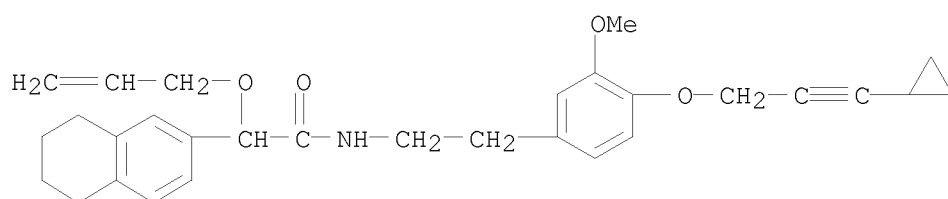
RN 1055230-44-8 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

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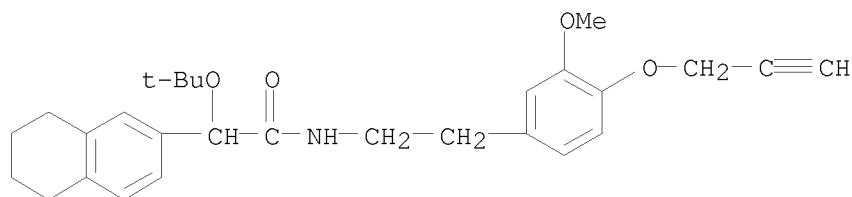
RN 1055230-47-1 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(2-propen-1-yloxy)- (CA INDEX NAME)



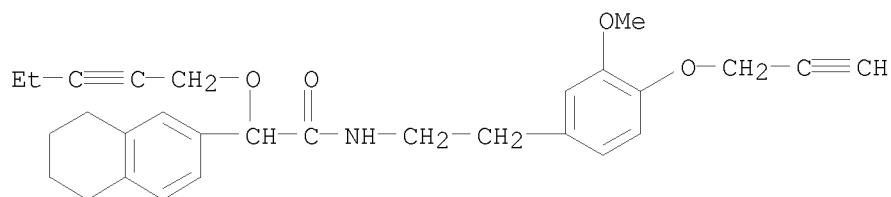
RN 1055230-72-2 CAPLUS

CN 2-Naphthaleneacetamide, α-(1,1-dimethylethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055230-99-3 CAPLUS

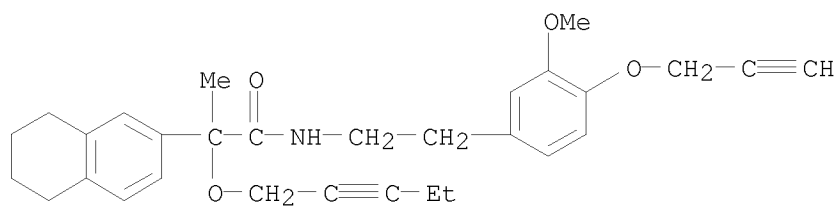
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(2-pentyn-1-yloxy)- (CA INDEX NAME)



RN 1055231-00-9 CAPLUS

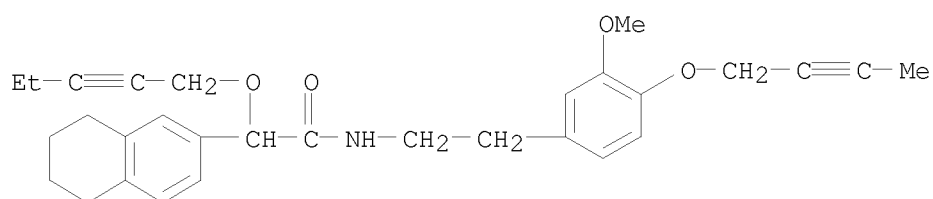
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-(2-pentyn-1-yloxy)- (CA INDEX NAME)

10/513699



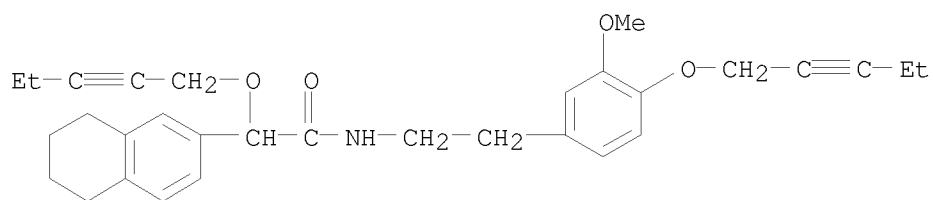
RN 1055231-01-0 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(2-pentyn-1-yloxy)- (CA INDEX NAME)



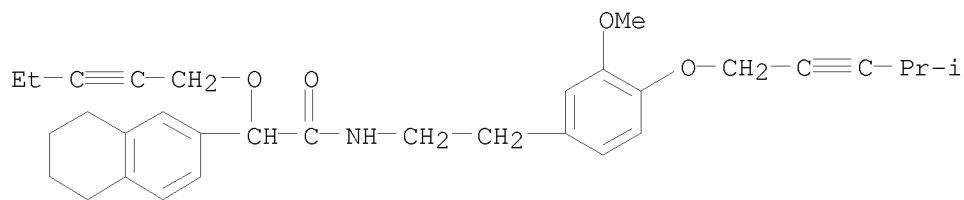
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CN INDEX NAME NOT YET ASSIGNED



RN 1055231-03-2 CAPLUS

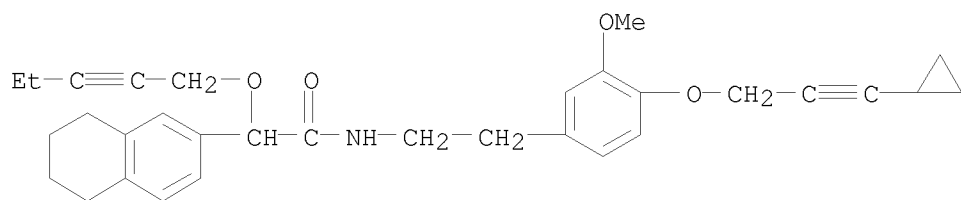
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RN 1055231-04-3 CAPLUS

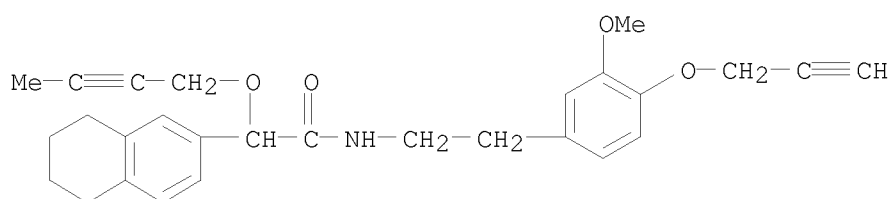
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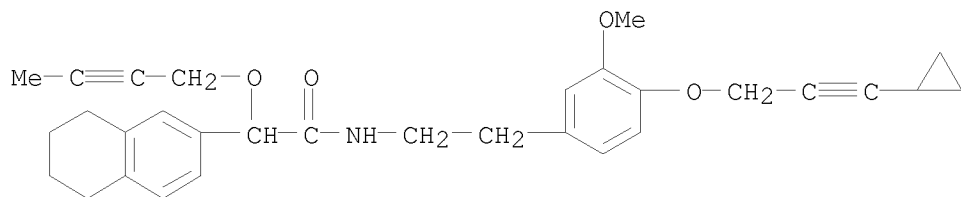
RN 1055232-70-6 CAPLUS

CN 2-Naphthaleneacetamide, α -(2-butyn-1-yloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



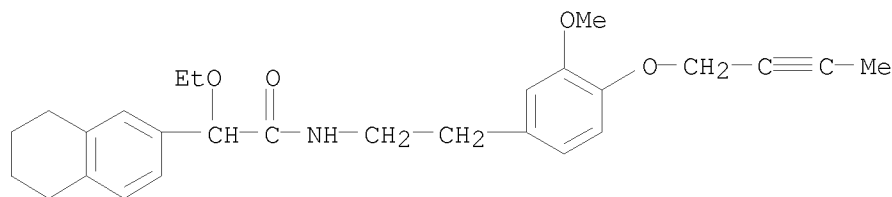
RN 1055232-71-7 CAPLUS

CN 2-Naphthaleneacetamide, α -(2-butyn-1-yloxy)-N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 1055236-78-6 CAPLUS

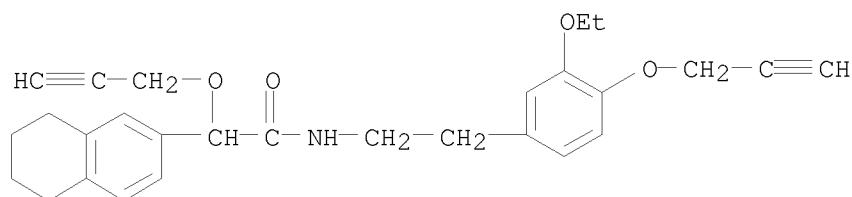
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -ethoxy-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 1055238-63-5 CAPLUS

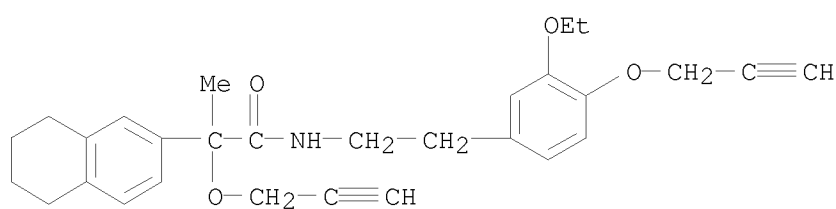
CN 2-Naphthaleneacetamide, N-[2-[3-ethoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)

10/513699



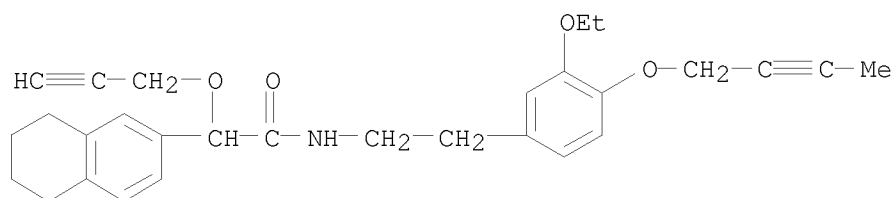
RN 1055238-64-6 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[3-ethoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-5,6,7,8-tetrahydro-α-methyl-α-(2-propyn-1-yloxy)- (CA INDEX NAME)



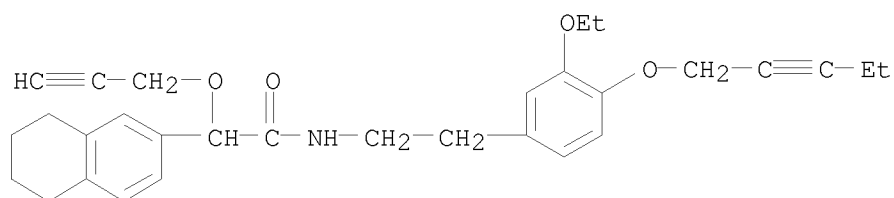
RN 1055238-65-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-ethoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055238-66-8 CAPLUS

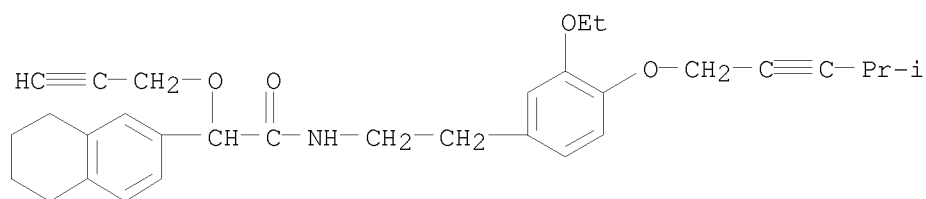
CN 2-Naphthaleneacetamide, N-[2-[3-ethoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]-5,6,7,8-tetrahydro-α-(2-propyn-1-yloxy)- (CA INDEX NAME)



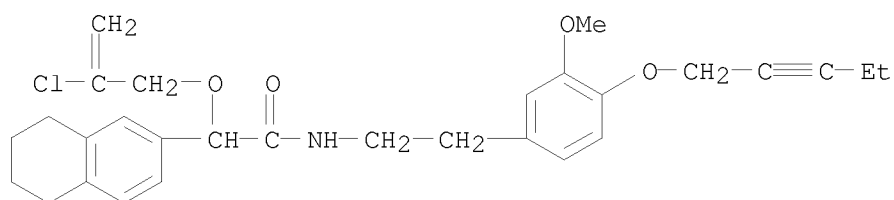
RN 1055238-67-9 CAPLUS

CN INDEX NAME NOT YET ASSIGNED

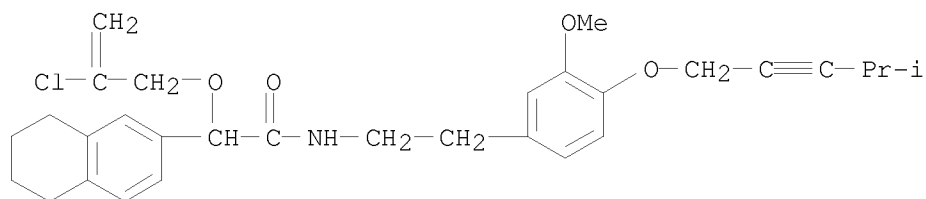
10/513699



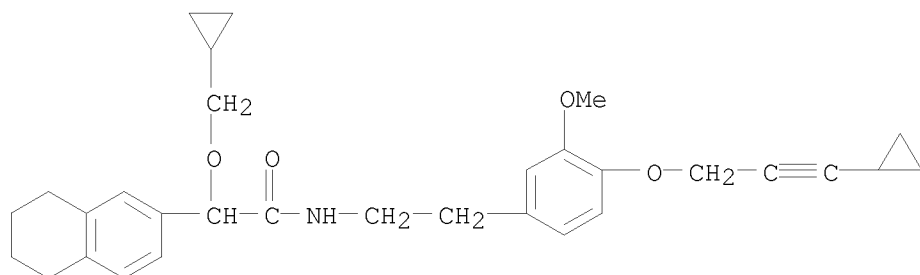
RN 1055240-59-9 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055240-60-2 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

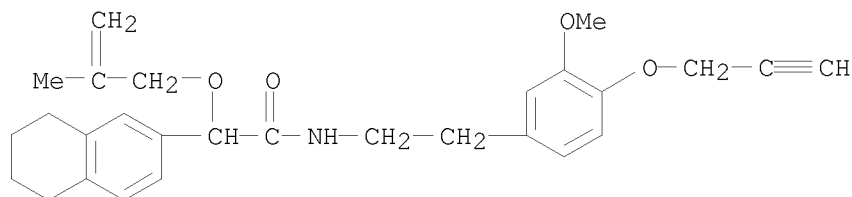


RN 1055242-23-3 CAPLUS
CN 2-Naphthaleneacetamide, α -(cyclopropylmethoxy)-N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)



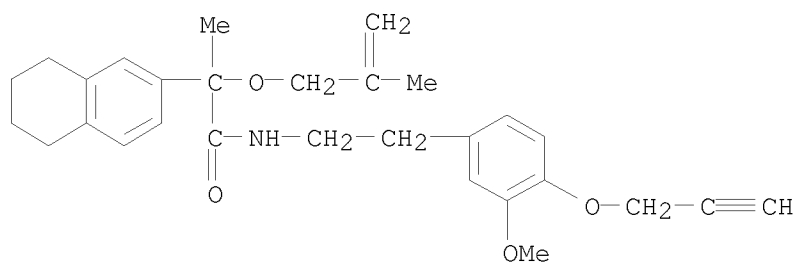
RN 1055243-85-0 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(2-methyl-2-propen-1-yl)oxy)- (CA INDEX NAME)

10/513699



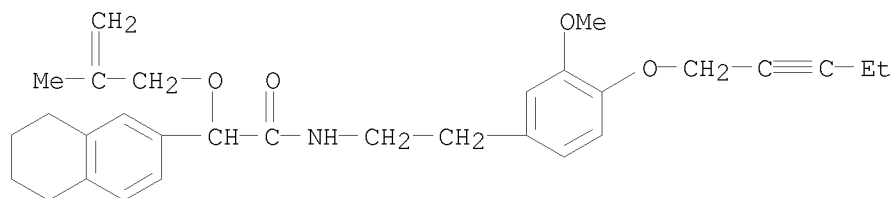
RN 1055243-86-1 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-[(2-methyl-2-propen-1-yl)oxy]-
(CA INDEX NAME)



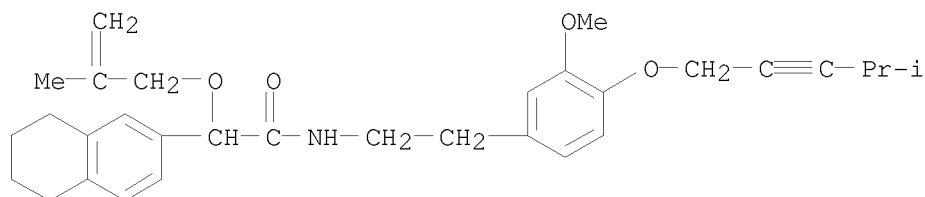
RN 1055243-87-2 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



RN 1055243-88-3 CAPLUS

CN INDEX NAME NOT YET ASSIGNED

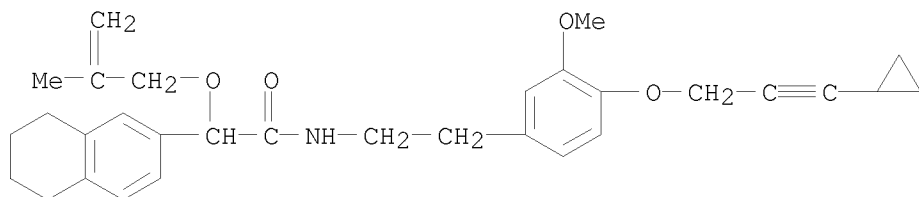


RN 1055243-89-4 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-[(2-methyl-2-propen-1-

10/513699

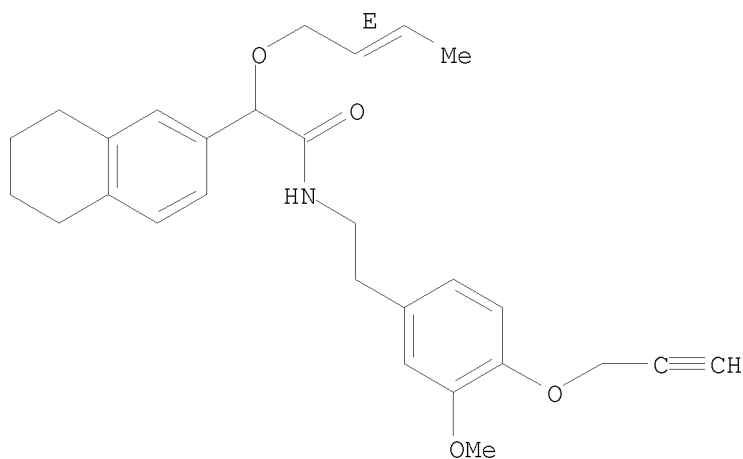
y1)oxy]- (CA INDEX NAME)



RN 1055244-11-5 CAPLUS

CN 2-Naphthaleneacetamide, α -[(2E)-2-buten-1-yloxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

Double bond geometry as shown.

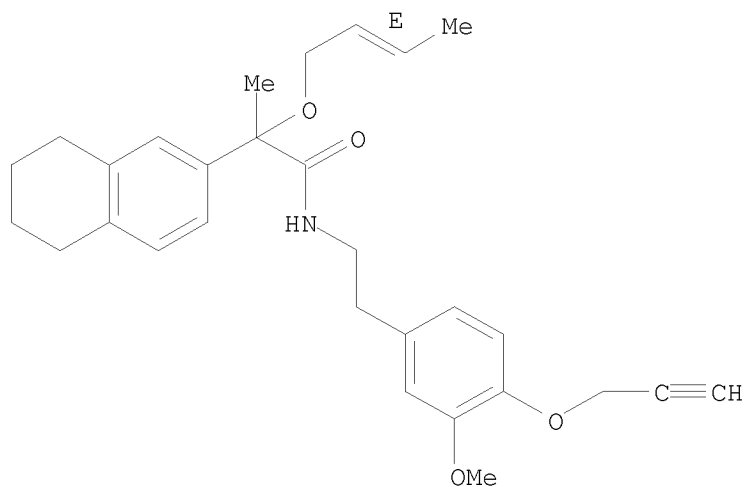


RN 1055244-12-6 CAPLUS

CN 2-Naphthaleneacetamide, α -[(2E)-2-buten-1-yloxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)

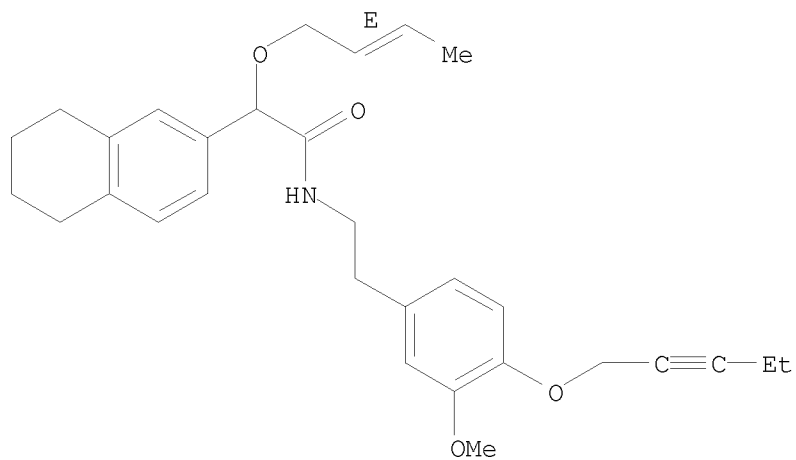
Double bond geometry as shown.

10/513699



RN 1055244-13-7 CAPLUS
CN 2-Naphthaleneacetamide, α -[(2E)-2-buten-1-yloxy]-5,6,7,8-tetrahydro-
N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

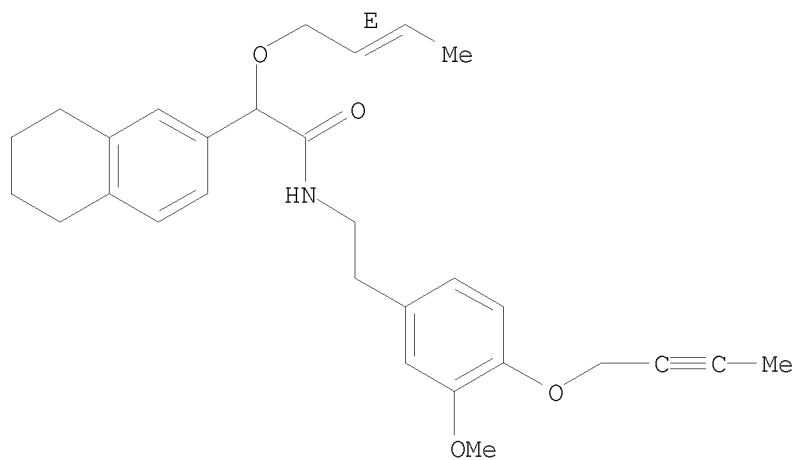
Double bond geometry as shown.



RN 1055244-14-8 CAPLUS
CN 2-Naphthaleneacetamide, α -[(2E)-2-buten-1-yloxy]-N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)

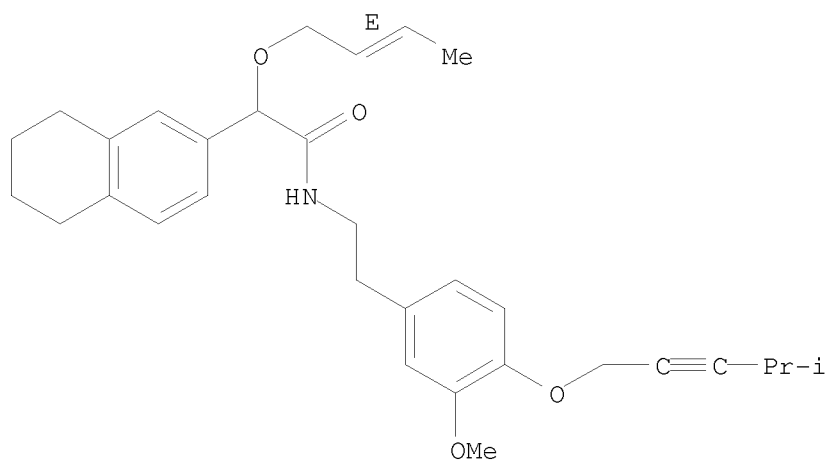
Double bond geometry as shown.

10/513699

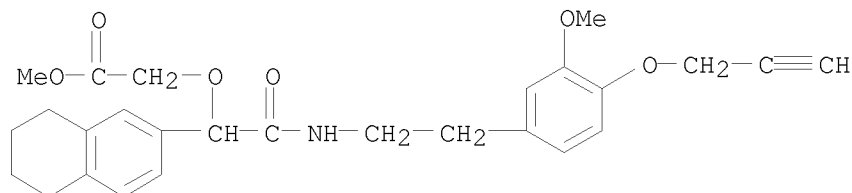


RN 1055244-15-9 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

Double bond geometry as shown.



RN 1055247-12-5 CAPLUS
CN Acetic acid, 2-[2-[[2-[[2-(3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-2-oxo-1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethoxy]-, methyl ester (CA INDEX NAME)



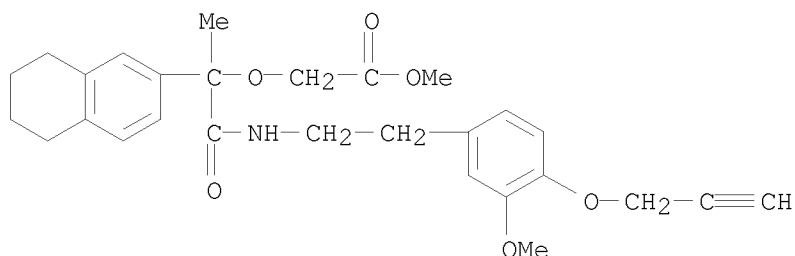
<12/04/2007>

Erich Leese

10/513699

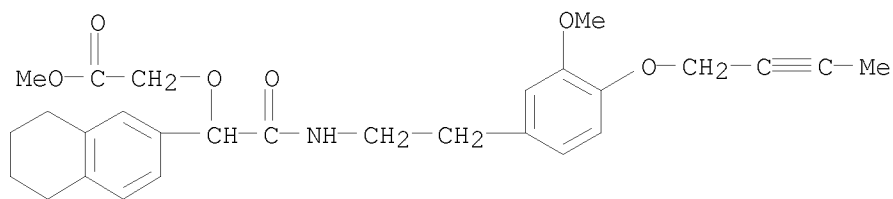
RN 1055247-15-8 CAPLUS

CN Acetic acid, 2-[2-[[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]amino]-1-methyl-2-oxo-1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethoxy]-, methyl ester (CA INDEX NAME)



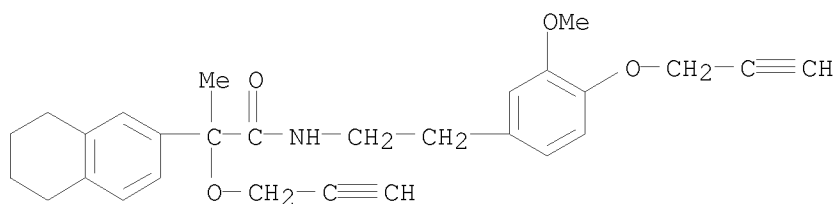
RN 1055247-16-9 CAPLUS

CN Acetic acid, 2-[2-[[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]amino]-2-oxo-1-(5,6,7,8-tetrahydro-2-naphthalenyl)ethoxy]-, methyl ester (CA INDEX NAME)



RN 1055249-28-9 CAPLUS

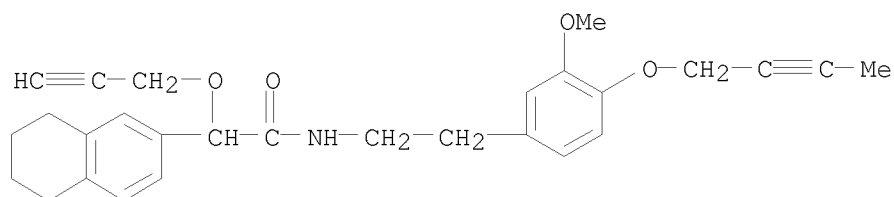
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055249-29-0 CAPLUS

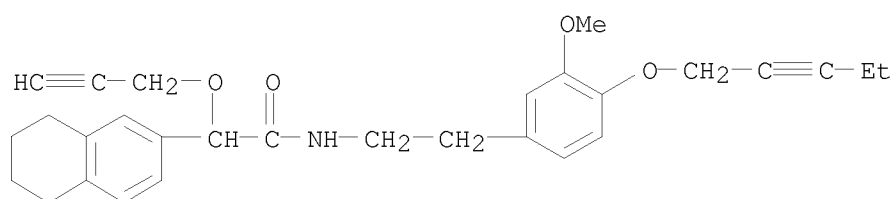
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)

10/513699



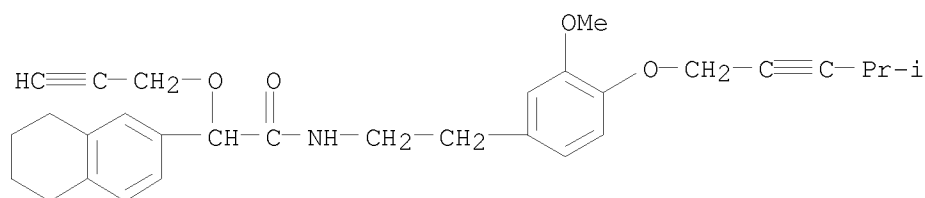
RN 1055249-30-3 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]-α-(2-propyn-1-yloxy)- (CA INDEX NAME)



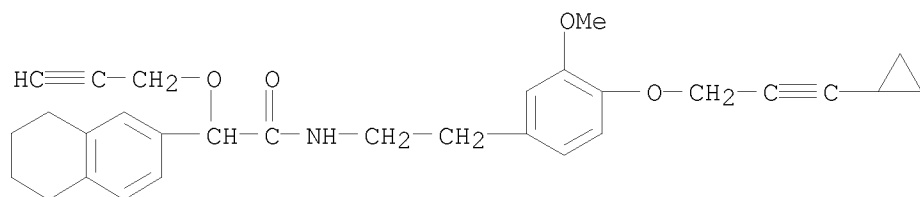
RN 1055249-33-6 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



RN 1055249-34-7 CAPLUS

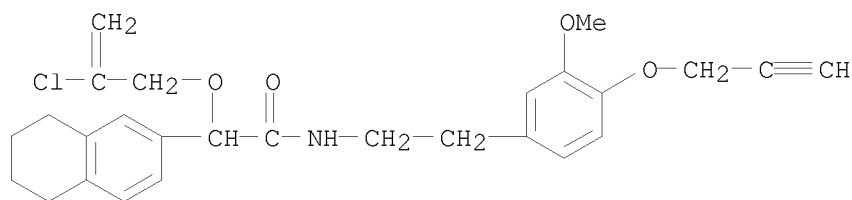
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1055250-39-9 CAPLUS

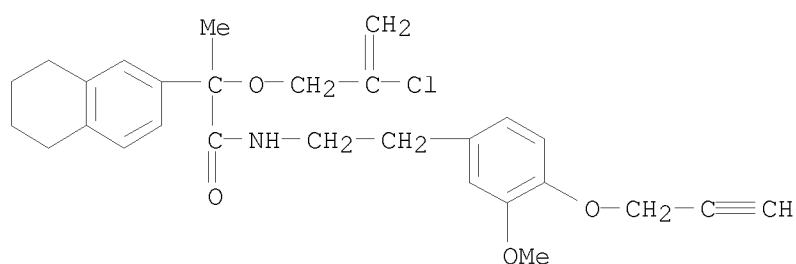
CN 2-Naphthaleneacetamide, α-[(2-chloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

10/513699



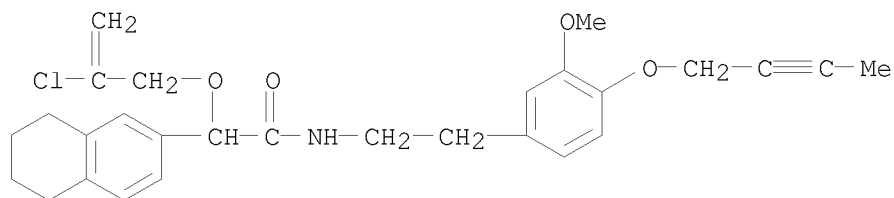
RN 1055250-41-3 CAPLUS

CN 2-Naphthaleneacetamide, α -[(2-chloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



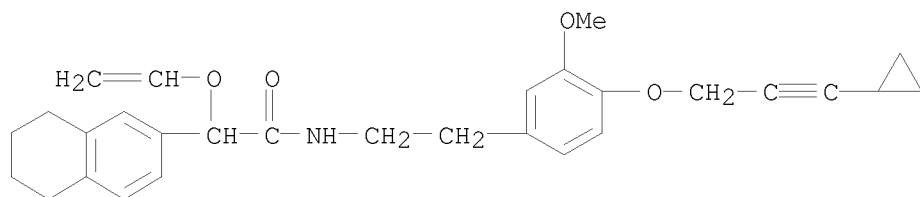
RN 1055250-45-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -[(2-chloro-2-propen-1-yl)oxy]-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 1055253-12-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]- α -(ethenyloxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)



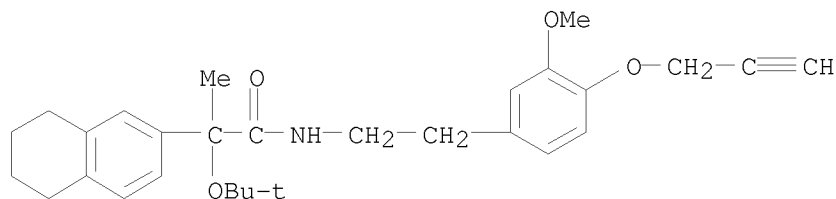
RN 1055254-60-8 CAPLUS

<12/04/2007>

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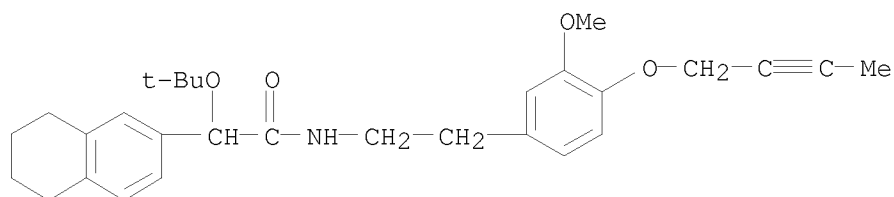
10/513699

CN 2-Naphthaleneacetamide, α -(1,1-dimethylethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



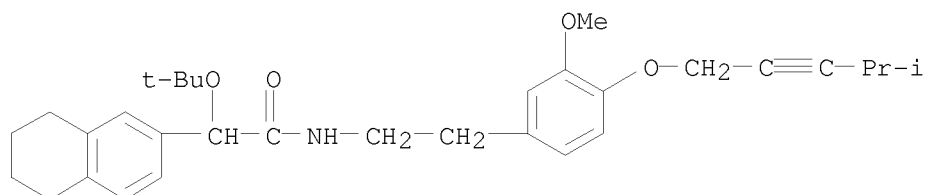
RN 1055254-61-9 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -(1,1-dimethylethoxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)



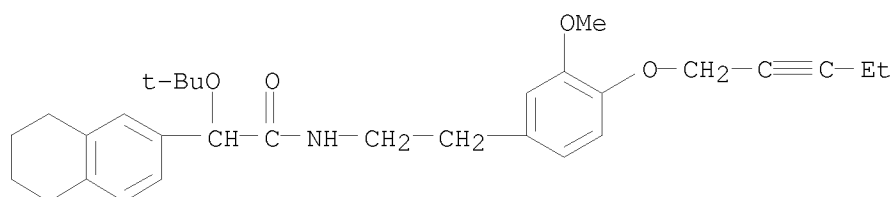
RN 1055254-62-0 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



RN 1055254-63-1 CAPLUS

CN 2-Naphthaleneacetamide, α -(1,1-dimethylethoxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

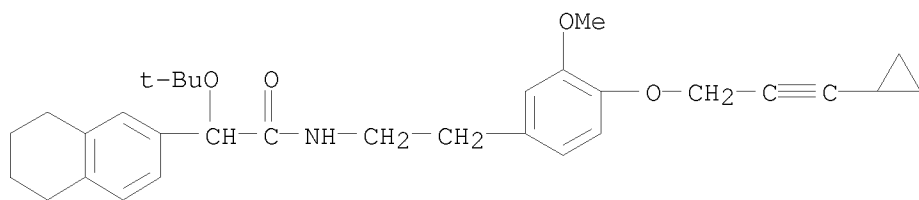


RN 1055254-64-2 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-

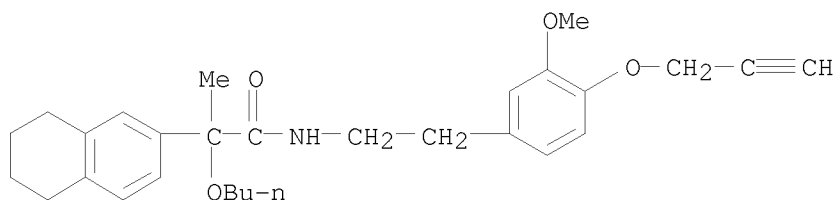
10/513699

methoxyphenyl]ethyl]- α -(1,1-dimethylethoxy)-5,6,7,8-tetrahydro- (CA INDEX NAME)



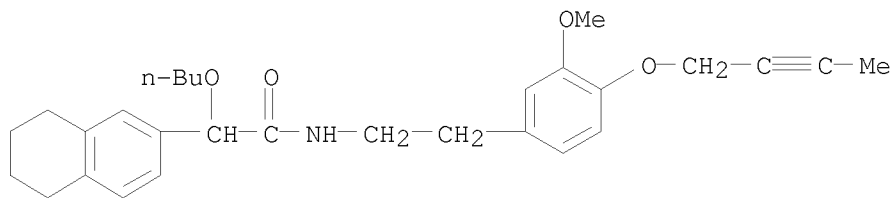
RN 1055255-69-0 CAPLUS

CN 2-Naphthaleneacetamide, α -butoxy-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



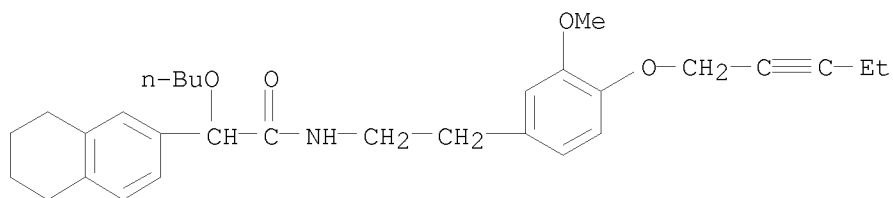
RN 1055255-70-3 CAPLUS

CN 2-Naphthaleneacetamide, α -butoxy-N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 1055255-71-4 CAPLUS

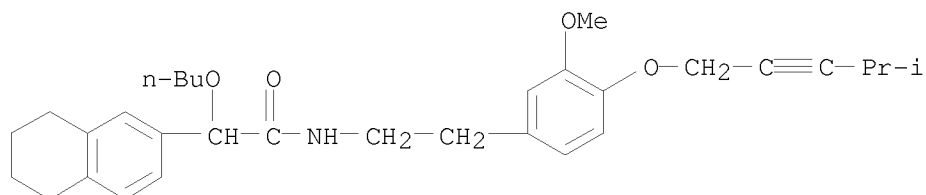
CN 2-Naphthaleneacetamide, α -butoxy-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055255-72-5 CAPLUS

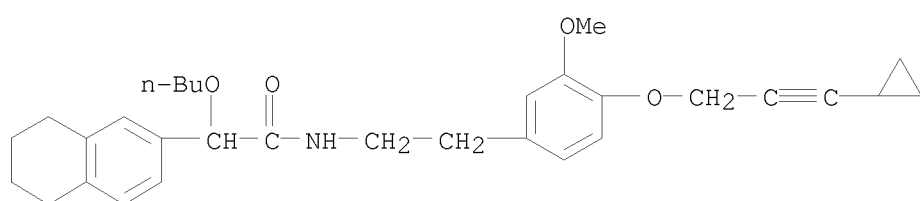
CN INDEX NAME NOT YET ASSIGNED

10/513699



RN 1055255-73-6 CAPLUS

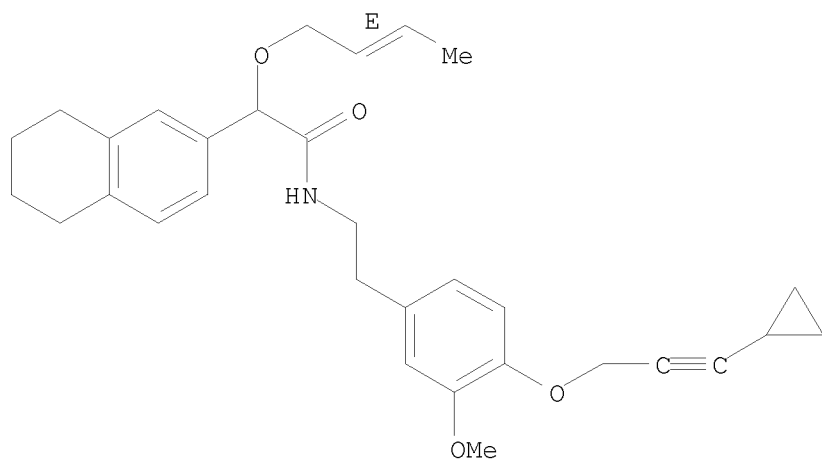
CN 2-Naphthaleneacetamide, α -butoxy-N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 1055256-73-9 CAPLUS

CN 2-Naphthaleneacetamide, α -[(2E)-2-buten-1-yloxy]-N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)

Double bond geometry as shown.



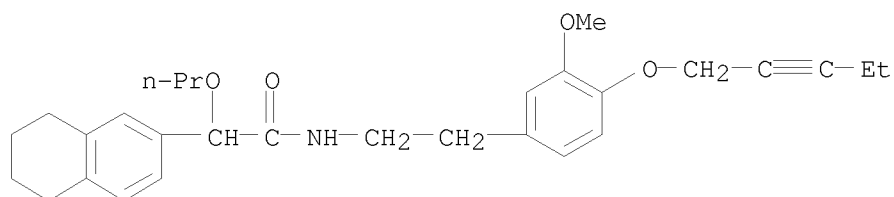
RN 1055258-31-5 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- α -propoxy- (CA INDEX NAME)

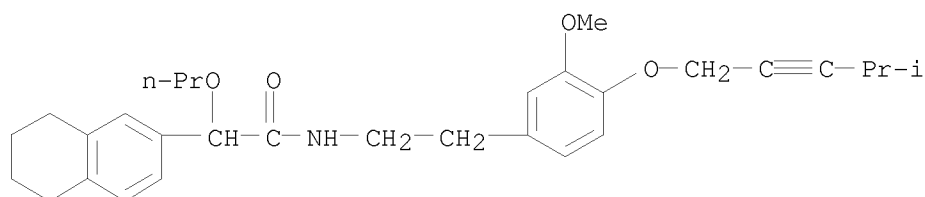
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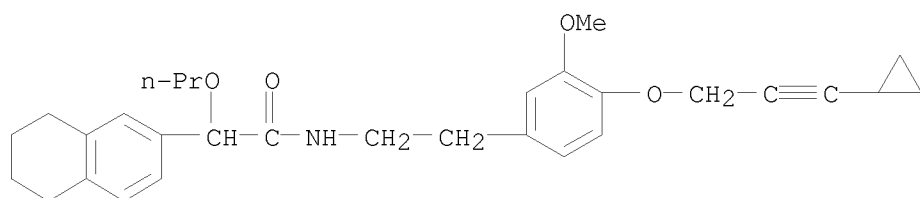
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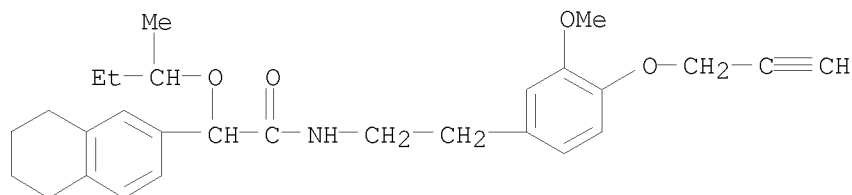
RN 1055258-32-6 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055258-33-7 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-propoxy- (CA INDEX NAME)

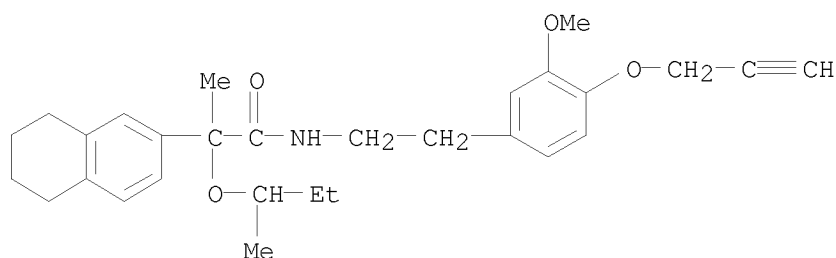


RN 1055261-20-5 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(1-methylpropoxy)- (CA INDEX NAME)



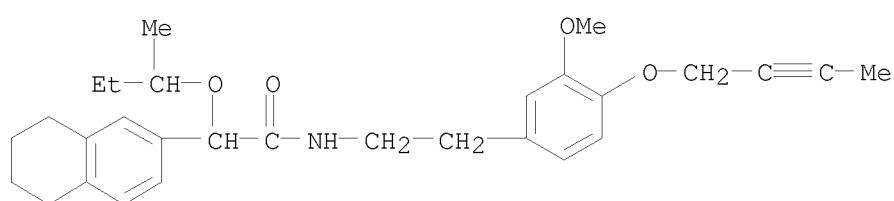
RN 1055261-21-6 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-(1-methylpropoxy)- (CA INDEX NAME)

10/513699



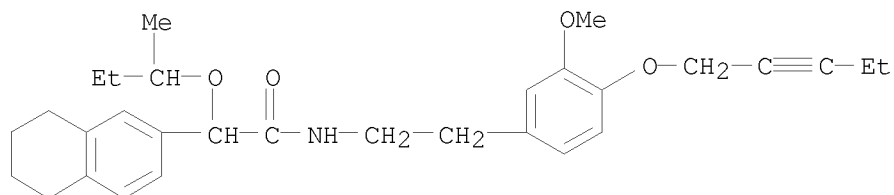
RN 1055261-22-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(1-methylpropoxy)- (CA INDEX NAME)



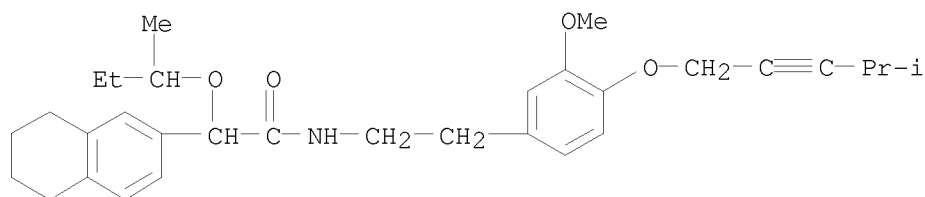
RN 1055261-23-8 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- α -(1-methylpropoxy)- (CA INDEX NAME)



RN 1055261-24-9 CAPLUS

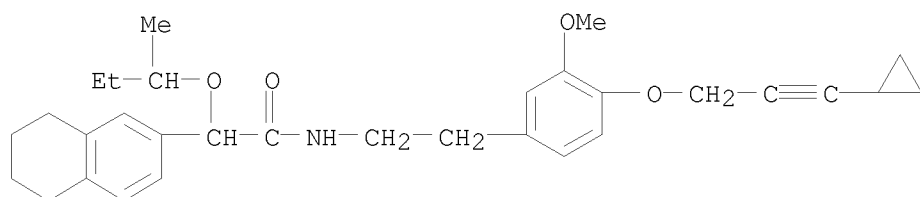
CN INDEX NAME NOT YET ASSIGNED



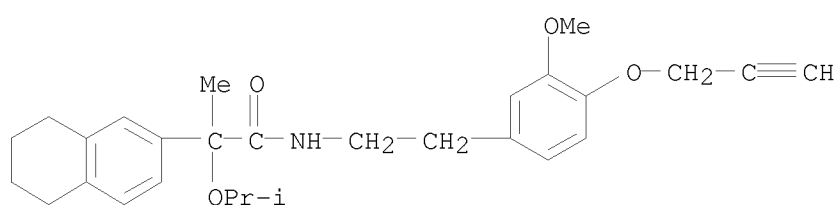
RN 1055261-25-0 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(1-methylpropoxy)- (CA INDEX NAME)

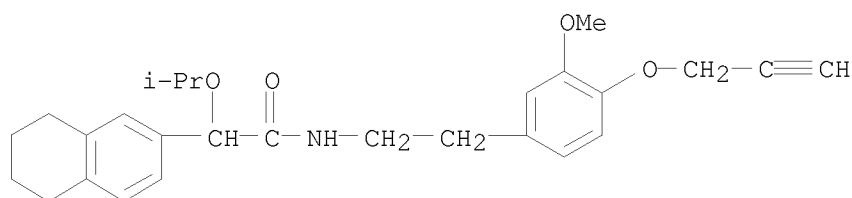
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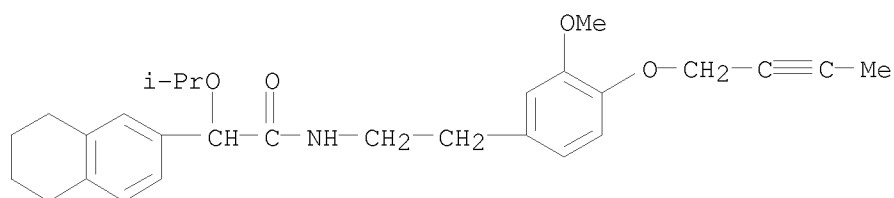
RN 1055262-68-4 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-(1-methylethoxy)- (CA INDEX NAME)



RN 1055262-70-8 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(1-methylethoxy)- (CA INDEX NAME)

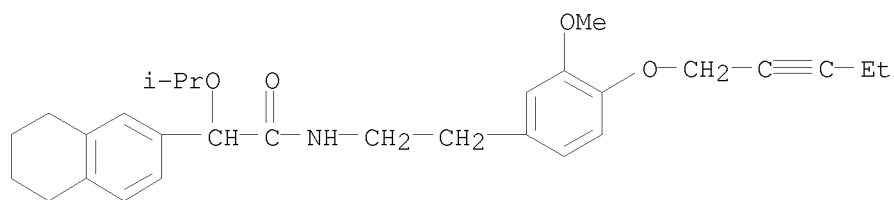


RN 1055262-73-1 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(1-methylethoxy)- (CA INDEX NAME)

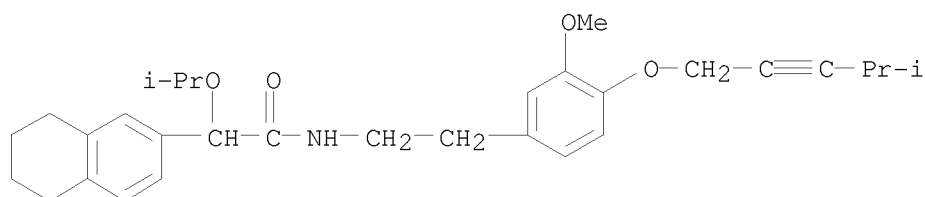


RN 1055262-75-3 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]-α-(1-methylethoxy)- (CA INDEX NAME)

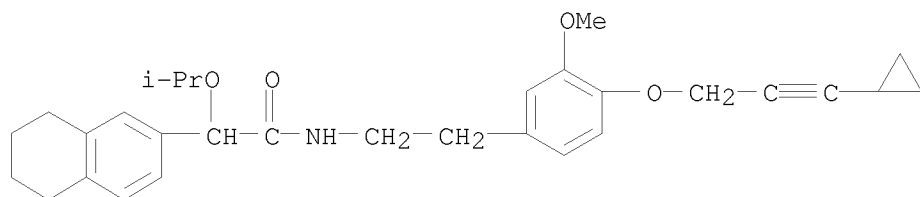
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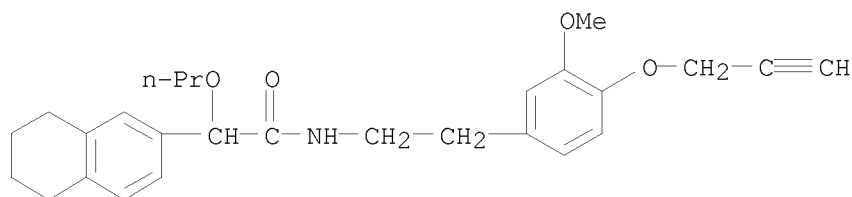
RN 1055262-76-4 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055262-78-6 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(1-methylethoxy)- (CA INDEX NAME)

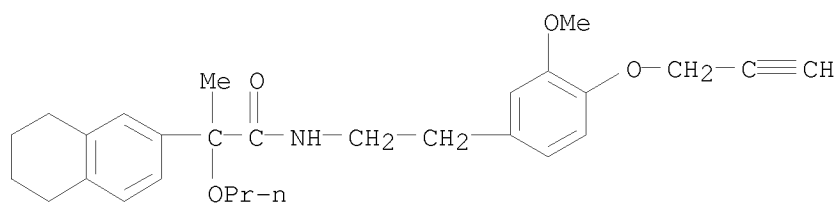


RN 1055264-42-0 CAPLUS
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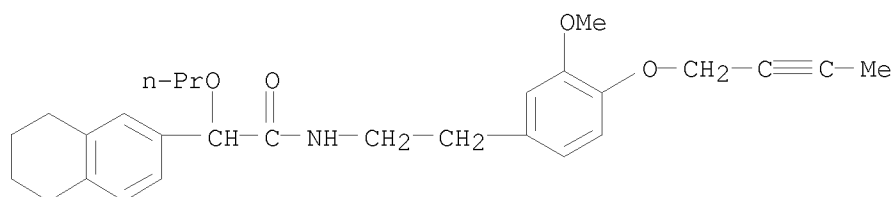
RN 1055264-43-1 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-propoxy- (CA INDEX NAME)

10/513699



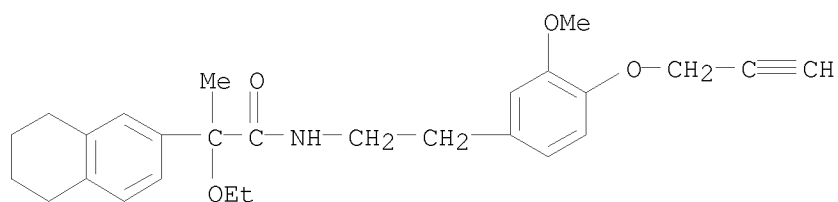
RN 1055264-45-3 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -propoxy- (CA INDEX NAME)



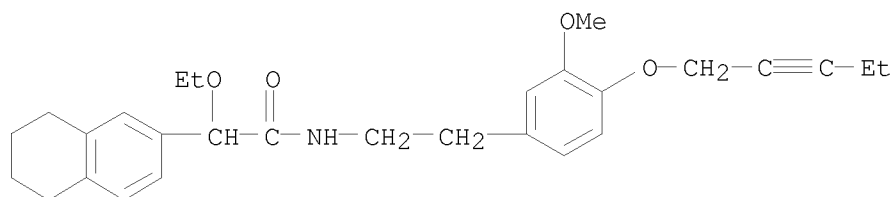
RN 1055267-02-1 CAPLUS

CN 2-Naphthaleneacetamide, α -ethoxy-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



RN 1055267-05-4 CAPLUS

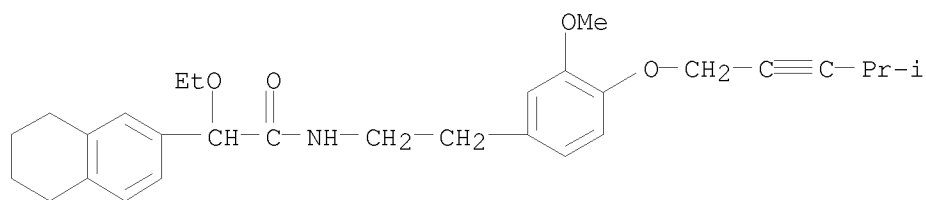
CN 2-Naphthaleneacetamide, α -ethoxy-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



RN 1055267-06-5 CAPLUS

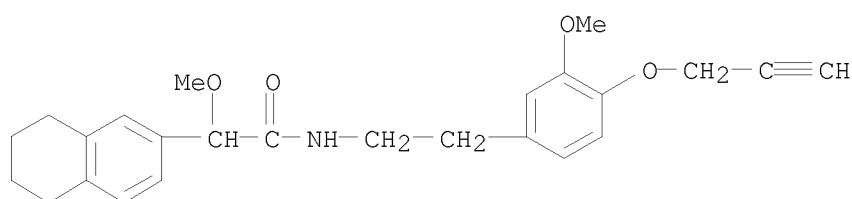
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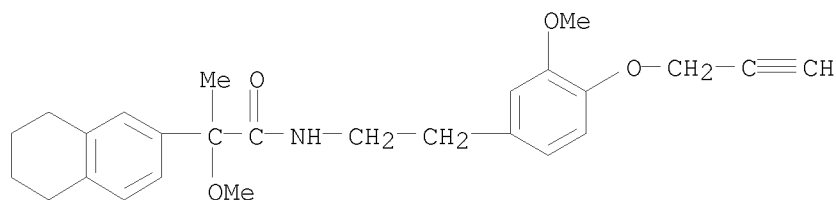
RN 1055270-60-4 CAPLUS

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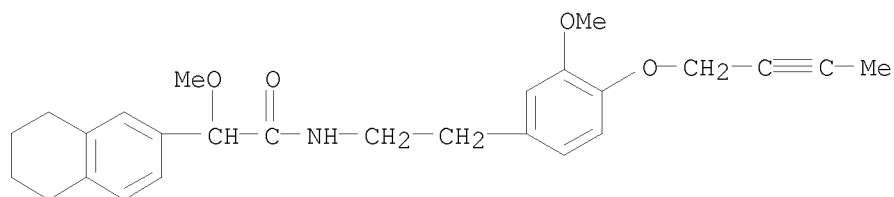
RN 1055270-61-5 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -methoxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



RN 1055270-62-6 CAPLUS

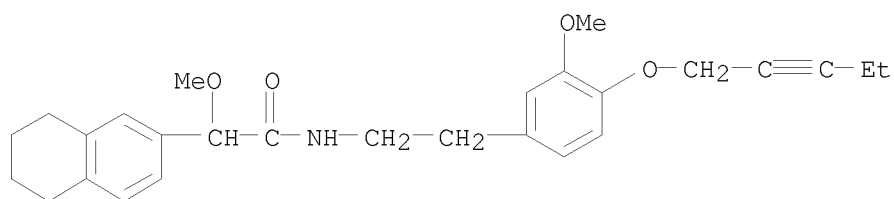
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -methoxy- (CA INDEX NAME)



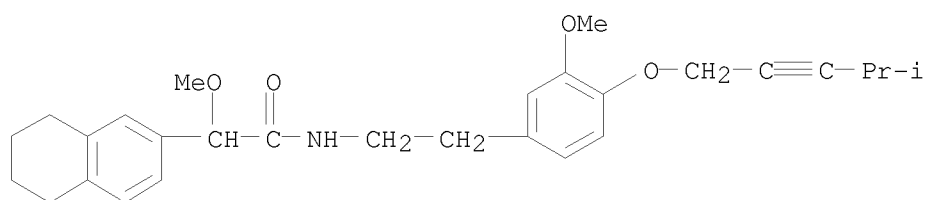
RN 1055270-63-7 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -methoxy-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

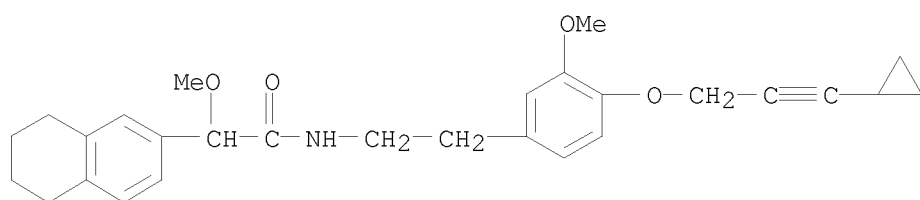
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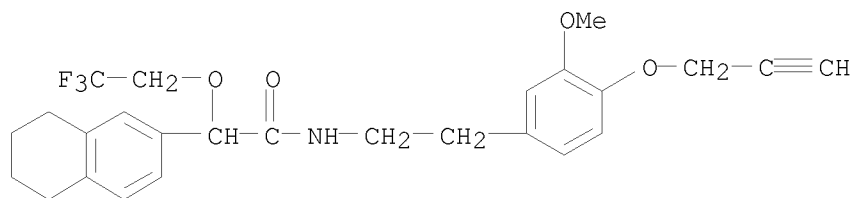
RN 1055270-64-8 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1055270-65-9 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-methoxy- (CA INDEX NAME)

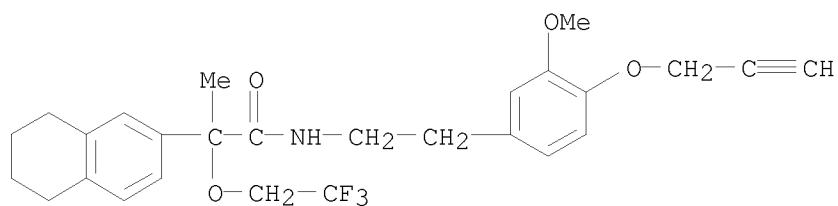


RN 1055271-92-5 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(2,2,2-trifluoroethoxy)- (CA INDEX NAME)



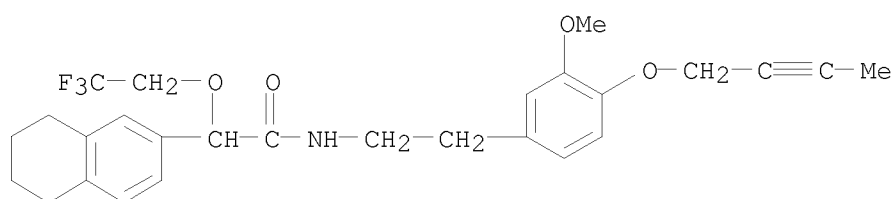
RN 1055271-93-6 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-(2,2,2-trifluoroethoxy)- (CA INDEX NAME)

10/513699



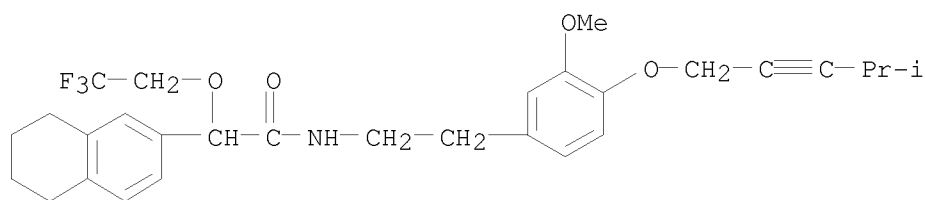
RN 1055271-94-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(2,2,2-trifluoroethoxy)- (CA INDEX NAME)



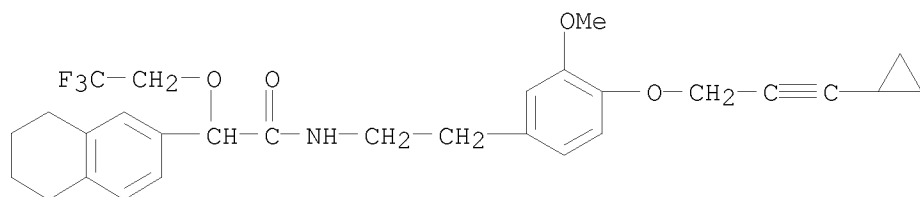
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CN INDEX NAME NOT YET ASSIGNED



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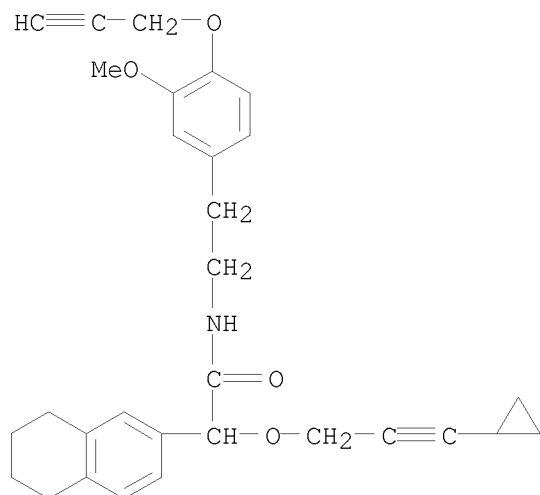
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(2,2,2-trifluoroethoxy)- (CA INDEX NAME)



RN 1055273-51-2 CAPLUS

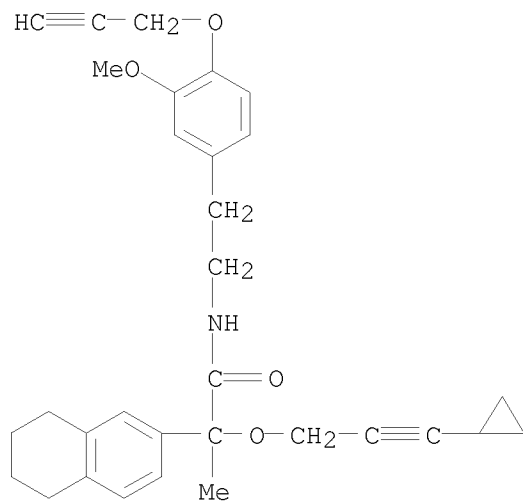
CN 2-Naphthaleneacetamide, α-[(3-cyclopropyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

10/513699



RN 1055273-52-3 CAPLUS

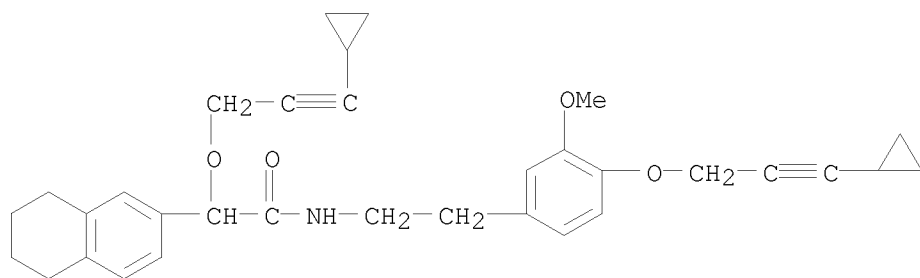
CN 2-Naphthaleneacetamide, α -[(3-cyclopropyl-2-propyn-1-yl)oxy]-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



RN 1055273-53-4 CAPLUS

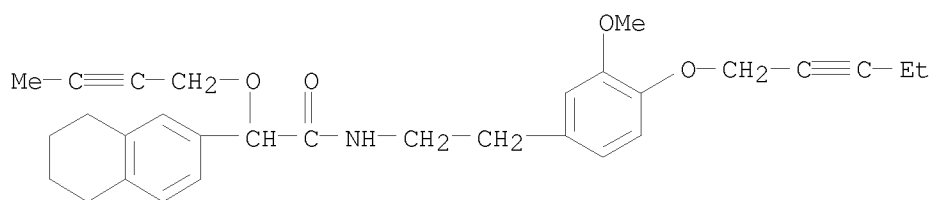
CN 2-Naphthaleneacetamide, α -[(3-cyclopropyl-2-propyn-1-yl)oxy]-N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)

10/513699



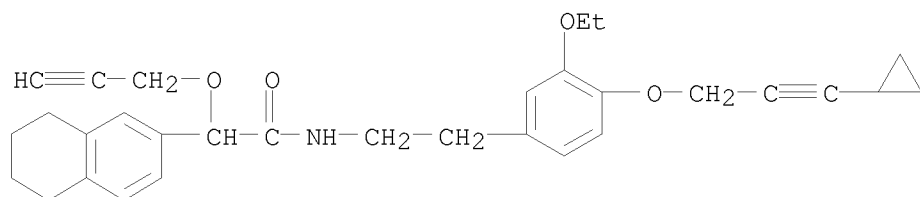
RN 1055274-16-2 CAPLUS

CN 2-Naphthaleneacetamide, α -(2-butyn-1-yloxy)-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



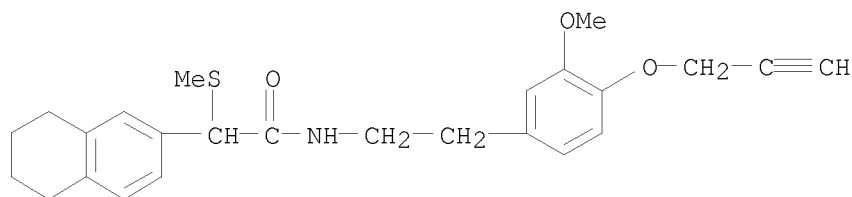
RN 1055275-96-1 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-ethoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(2-propyn-1-yloxy)- (CA INDEX NAME)



RN 1102336-75-3 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(methylthio)- (CA INDEX NAME)



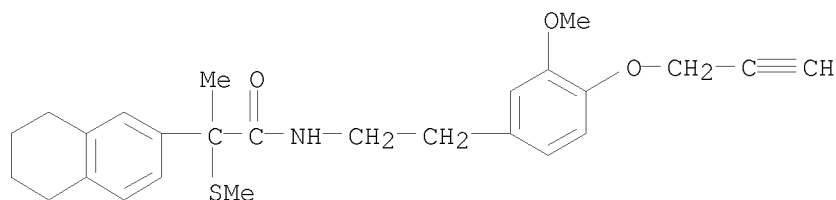
RN 1102336-76-4 CAPLUS

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Erich Leese

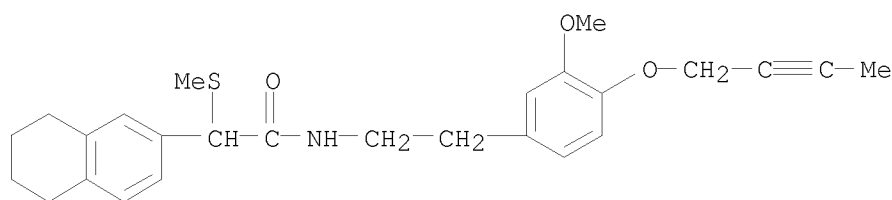
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CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- α -(methylthio)- (CA INDEX NAME)



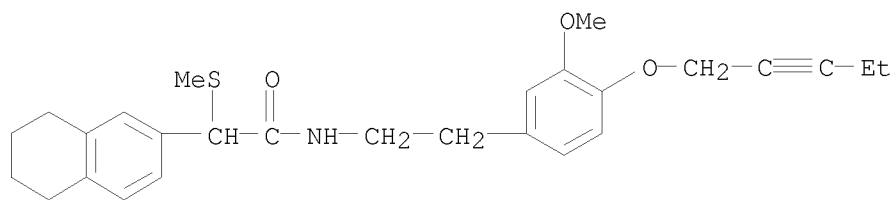
RN 1102336-77-5 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(methylthio)- (CA INDEX NAME)



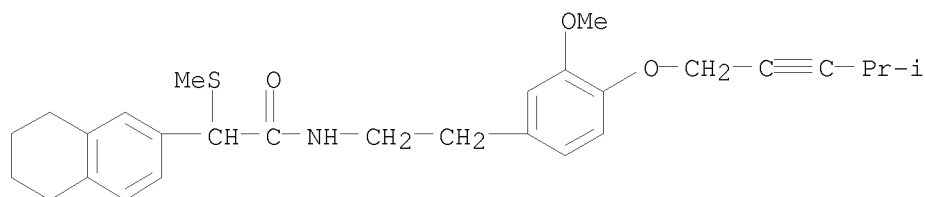
RN 1102336-78-6 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- α -(methylthio)- (CA INDEX NAME)



RN 1102336-79-7 CAPLUS

CN INDEX NAME NOT YET ASSIGNED

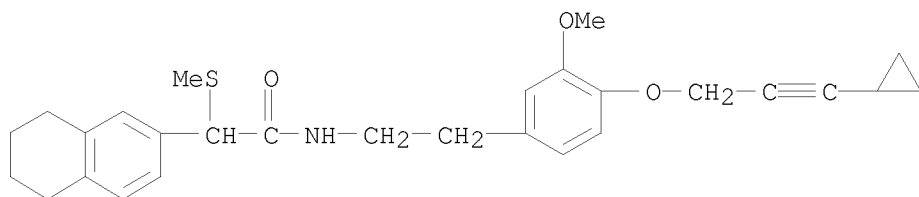


RN 1102336-80-0 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -(methylthio)- (CA INDEX NAME)

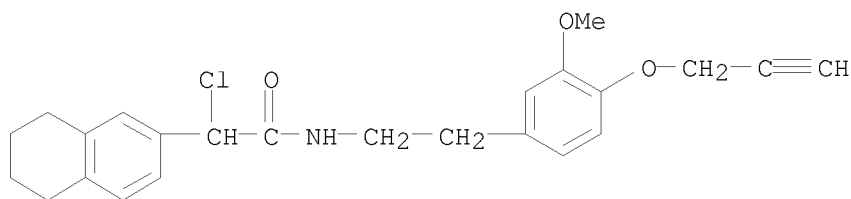
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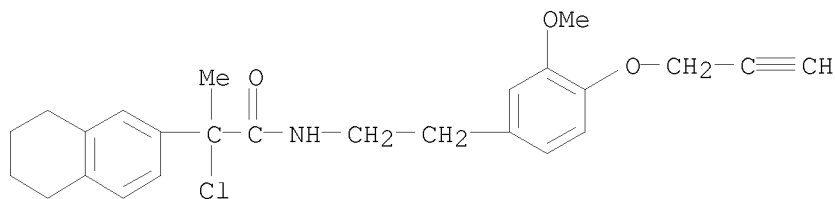
RN 1102339-25-2 CAPLUS

CN 2-Naphthaleneacetamide, α -chloro-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



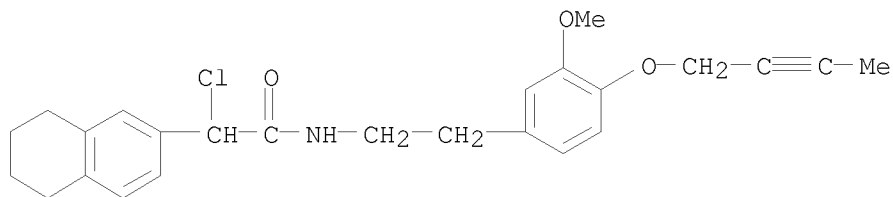
RN 1102339-26-3 CAPLUS

CN 2-Naphthaleneacetamide, α -chloro-5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- (CA INDEX NAME)



RN 1102339-27-4 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]- α -chloro-5,6,7,8-tetrahydro- (CA INDEX NAME)



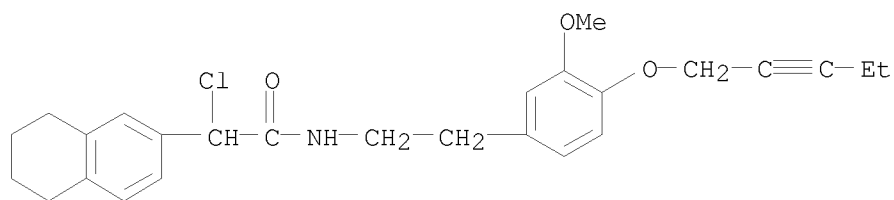
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CN INDEX NAME NOT YET ASSIGNED

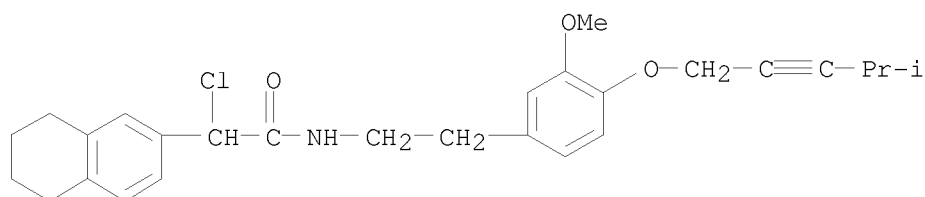
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Erich Leese

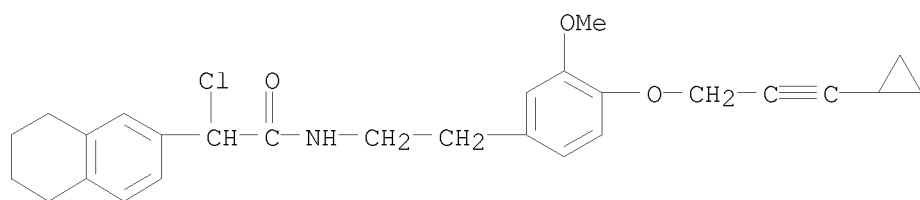
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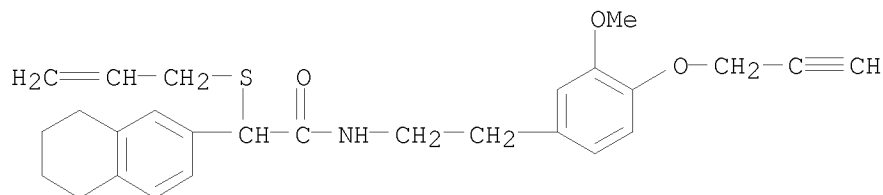
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CN INDEX NAME NOT YET ASSIGNED



RN 1102339-30-9 CAPLUS
CN 2-Naphthaleneacetamide, α -chloro-N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- (CA INDEX NAME)

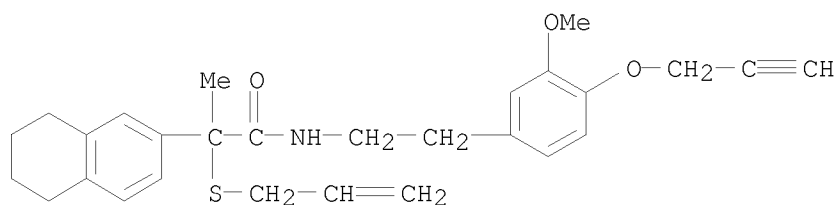


RN 1102340-45-3 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -(2-propen-1-ylthio)- (CA INDEX NAME)



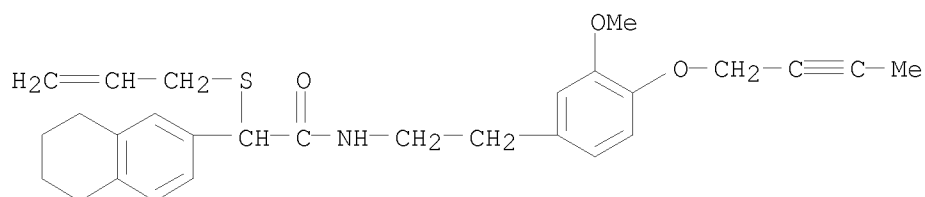
RN 1102340-46-4 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- α -methyl- α -(2-propen-1-ylthio)- (CA INDEX NAME)

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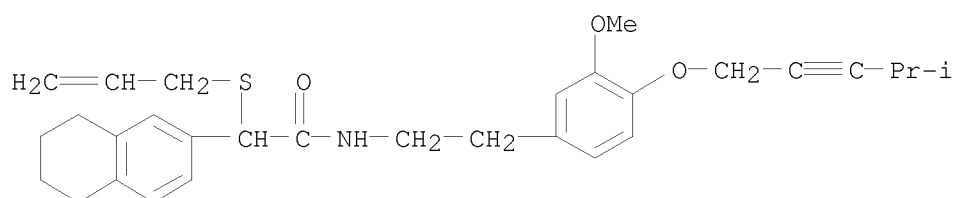
RN 1102340-47-5 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(2-propen-1-ylthio)- (CA INDEX NAME)



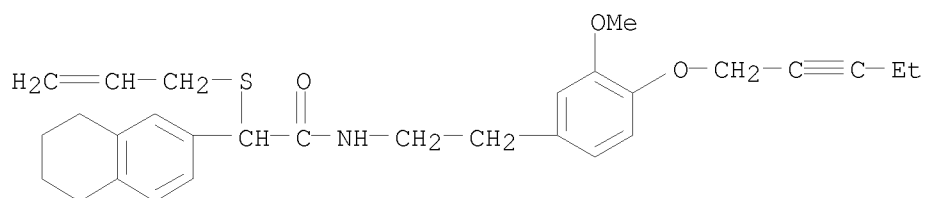
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CN INDEX NAME NOT YET ASSIGNED



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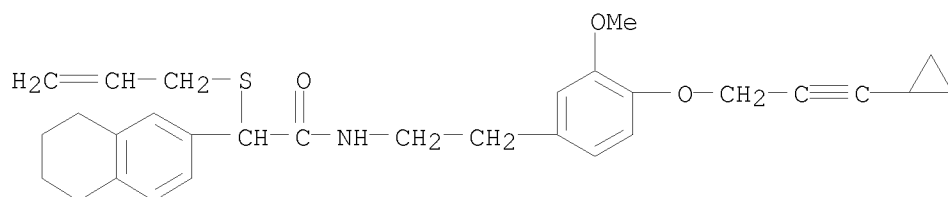
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]-α-(2-propen-1-ylthio)- (CA INDEX NAME)



RN 1102340-50-0 CAPLUS

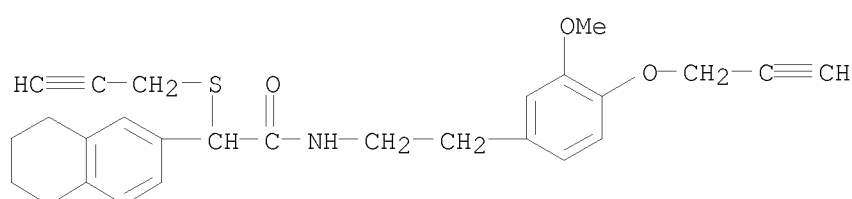
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(2-propen-1-ylthio)- (CA INDEX NAME)

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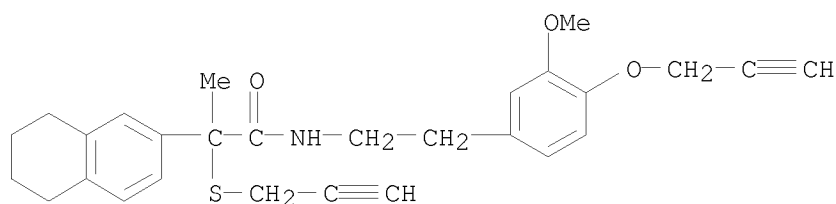
RN 1102343-22-5 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-(2-propyn-1-ylthio)- (CA INDEX NAME)



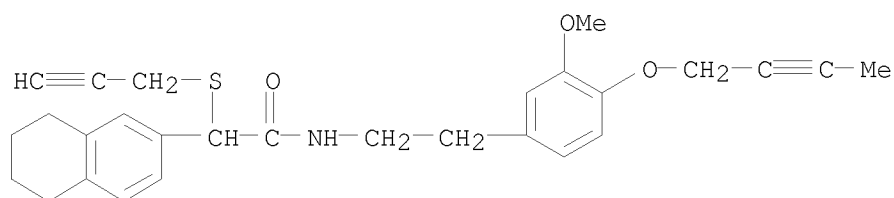
RN 1102343-23-6 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-methyl-α-(2-propyn-1-ylthio)- (CA INDEX NAME)



RN 1102343-24-7 CAPLUS

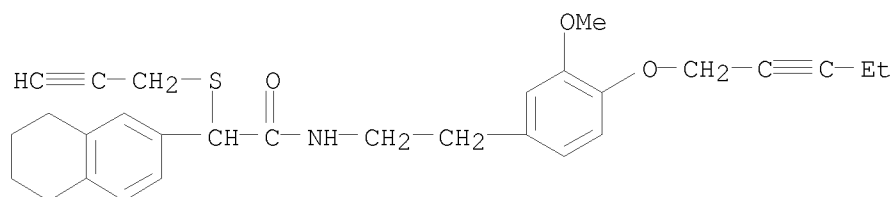
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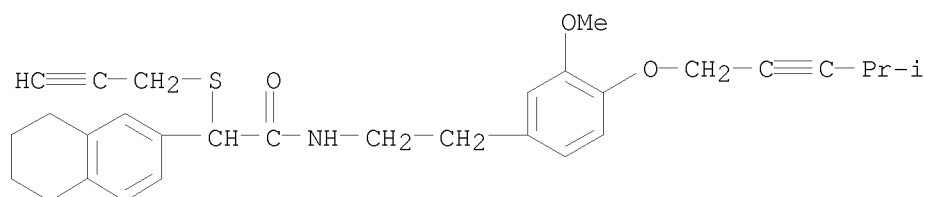
RN 1102343-25-8 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]-α-(2-propyn-1-ylthio)- (CA INDEX NAME)

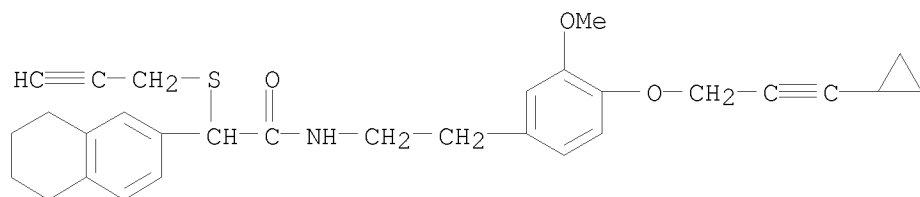
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RN 1102343-26-9 CAPLUS
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RN 1102343-27-0 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-(2-propyn-1-ylthio)- (CA INDEX NAME)



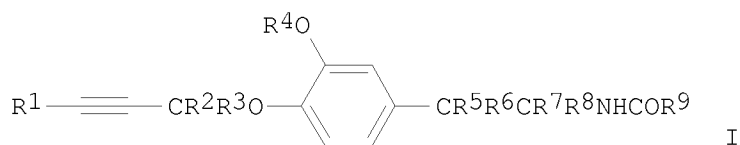
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REFERENCE COUNT: 6 THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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 ACCESSION NUMBER: 2000:493502 CAPLUS
 DOCUMENT NUMBER: 133:104883
 TITLE: Preparation of (acylaminoethyl)aryl propargyl ethers
 as agrochemical microbicides.
 INVENTOR(S): Zeller, Martin; Jeanguenat, Andre; Lamberth, Clemens;
 Kunz, Walter
 PATENT ASSIGNEE(S): Novartis A.-G., Switz.; Novartis-Erfindungen
 SOURCE: PCT Int. Appl., 89 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2000041998	A1	20000720	WO 2000-EP106	20000110 <--
W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW				
RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
TW 564244	B	20031201	TW 1999-88121433	19991204 <--
CA 2356121	A1	20000720	CA 2000-2356121	20000110 <--
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EP 1140799	A1	20011010	EP 2000-901518	20000110 <--
EP 1140799	B1	20040102		
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BR 2000007469	A	20011030	BR 2000-7469	20000110 <--
TR 200102004	T2	20011221	TR 2001-2004	20000110 <--
HU 2001005039	A2	20020429	HU 2001-5039	20000110 <--
HU 2001005039	A3	20020528		
JP 2002534494	T	20021015	JP 2000-593567	20000110 <--
AU 759247	B2	20030410	AU 2000-22886	20000110 <--
AT 257148	T	20040115	AT 2000-901518	20000110 <--
ES 2213565	T3	20040901	ES 2000-901518	20000110 <--
RU 2237058	C2	20040927	RU 2001-121196	20000110 <--
CN 1257153	C	20060524	CN 2000-802706	20000110
IL 144105	A	20061231	IL 2000-144105	20000110
PL 200004	B1	20081128	PL 2000-348923	20000110
EG 23070	A	20040229	EG 2000-18	20000111 <--
ZA 2001005514	A	20021004	ZA 2001-5514	20010704 <--
IN 2001CN00951	A	20050304	IN 2001-CN951	20010706
MX 2001006978	A	20011011	MX 2001-6978	20010709 <--
US 6469005	B1	20021022	US 2001-903651	20010711 <--
PRIORITY APPLN. INFO.:			GB 1999-455	A 19990111
			WO 2000-EP106	W 20000110

ASSIGNMENT HISTORY FOR US PATENT AVAILABLE IN LSUS DISPLAY FORMAT

OTHER SOURCE(S): MARPAT 133:104883
 GI



AB Title compds. [I; R1 = H, alkyl, cycloalkyl, (substituted) aryl; R2, R3 = H, alkyl; R4 = alkyl, alkenyl, alkynyl; R5-R8 = H, alkyl; R9 = COR10, C(OZ)R10R11, C(:NOR12)R10, etc.; R10 = (substituted) aryl, heteroaryl; R11 = H (substituted) alkyl, alkenyl, alkynyl; Z = H, COR16, CO2R16, COCO2R16, CONR16R17; R12 = H, (substituted) alkyl, alkenyl, alkynyl; R16, R17 = H, (substituted) alkyl, cycloalkyl, aryl, heteroaryl], were prepared Thus, 2-(3,4-dichlorophenyl)-N-[2-(3-methoxy-4-prop-2-ynyloxyphenyl)ethyl]-2-oxoacetamide was heated with MeONH2.HCl and pyridine in EtOH at 80° for 4 h to give 2-(3,4-dichlorophenyl)-2-methoxyimino-N-[2-(3-methoxy-4-prop-2-ynyloxyphenyl)ethyl]acetamide. Numerous I as 0.02% sprays gave complete control of *Plasmopara viticola* on vines.

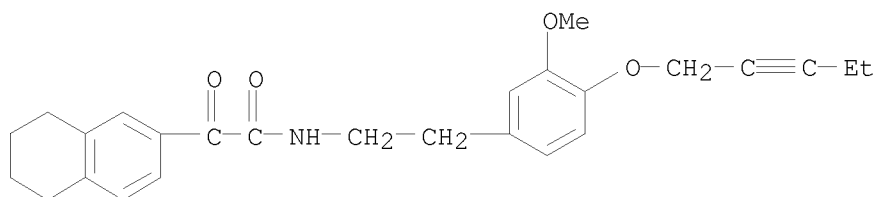
IT	1100606-32-3	1100607-98-4	1100609-21-9
	1100610-64-7	1100612-05-2	1100613-48-6
	1100614-54-7	1100616-02-1	1100617-03-5
	1100618-95-8	1100619-63-3	1100620-98-1
	1100621-30-4	1100623-97-9	1100626-75-2
	1100628-19-0	1100629-73-9	1100630-98-5
	1100633-66-6	1100634-28-3	

RL: PRPH (Prophetic)

(Preparation of (acylaminoethyl)aryl propargyl ethers as agrochemical microbicides.)

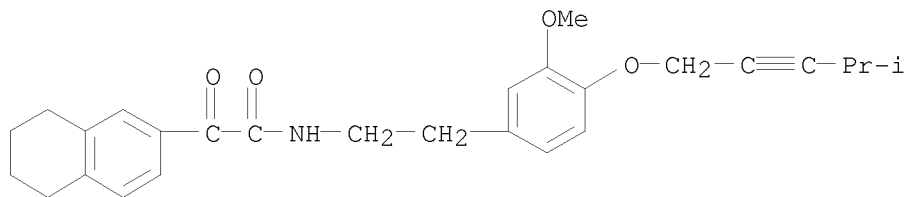
RN 1100606-32-3 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- α -oxo- (CA INDEX NAME)



RN 1100607-98-4 CAPLUS

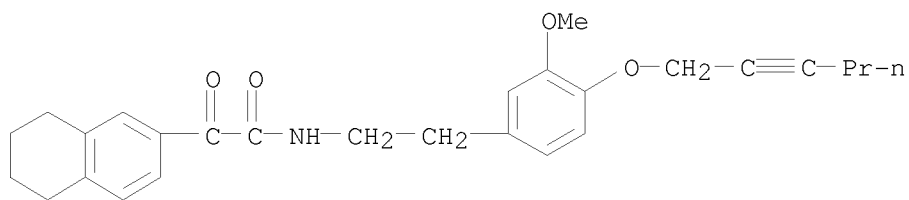
CN INDEX NAME NOT YET ASSIGNED



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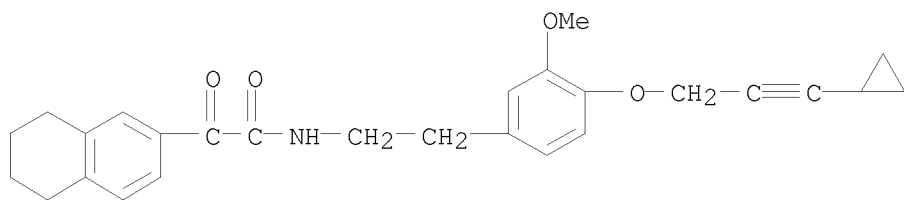
10/513699

CN 2-Naphthaleneacetamide, N-[2-[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -oxo- (CA INDEX NAME)



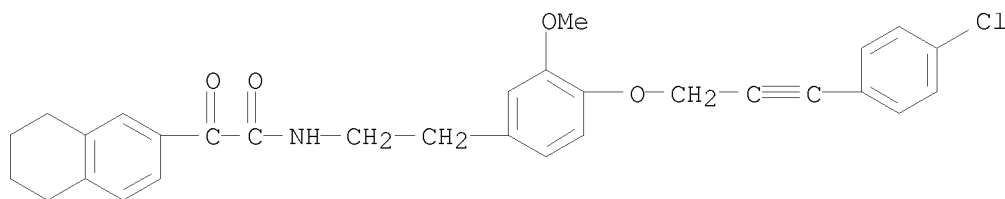
RN 1100610-64-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -oxo- (CA INDEX NAME)



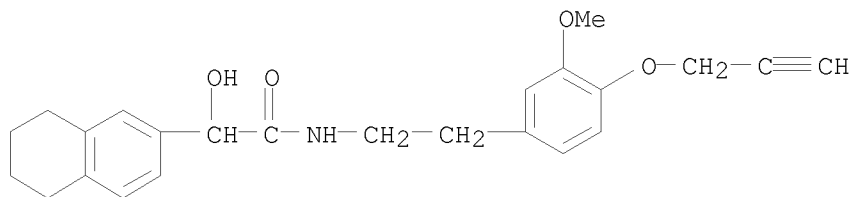
RN 1100612-05-2 CAPLUS

CN INDEX NAME NOT YET ASSIGNED



RN 1100613-48-6 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -hydroxy-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)



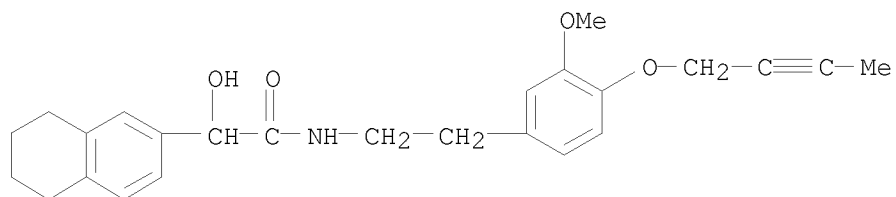
RN 1100614-54-7 CAPLUS

CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro- α -hydroxy- (CA INDEX NAME)

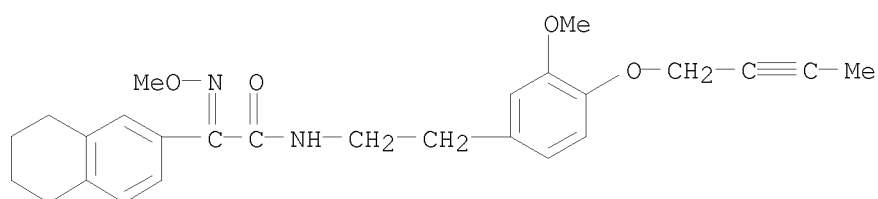
<12/04/2007>

Erich Leese

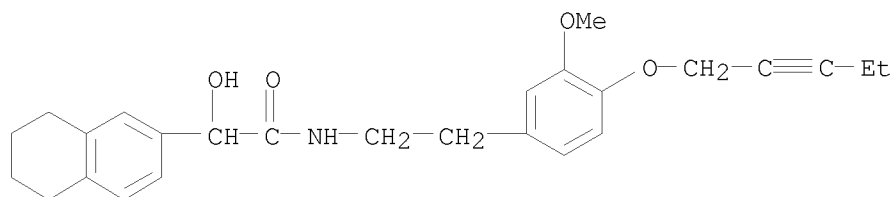
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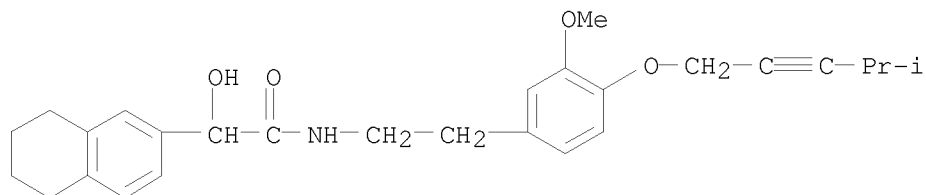
RN 1100616-02-1 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1100617-03-5 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro- α -hydroxy-N-[2-[3-methoxy-4-(2-pentyn-1-yloxy)phenyl]ethyl]- (CA INDEX NAME)

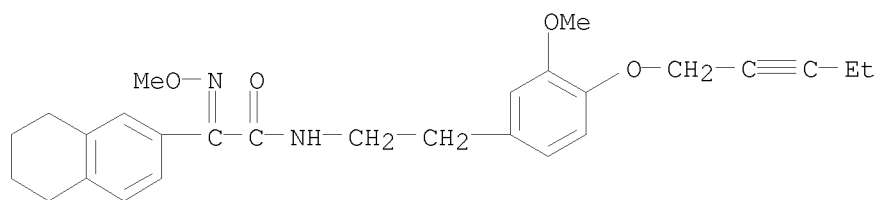


RN 1100618-95-8 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

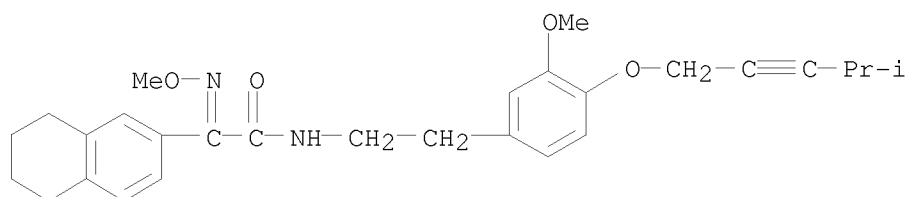


RN 1100619-63-3 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

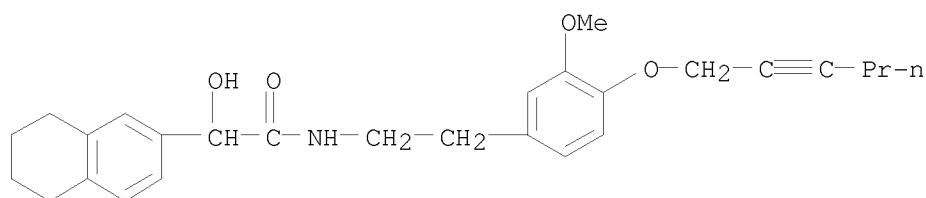
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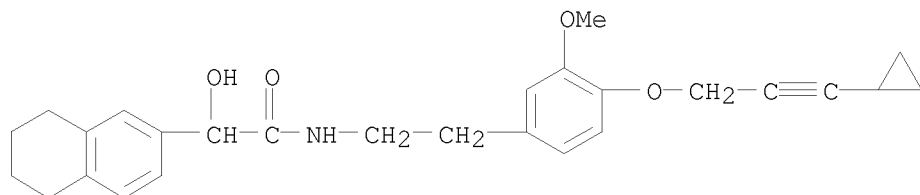
RN 1100620-98-1 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1100621-30-4 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-(2-hexyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-hydroxy- (CA INDEX NAME)

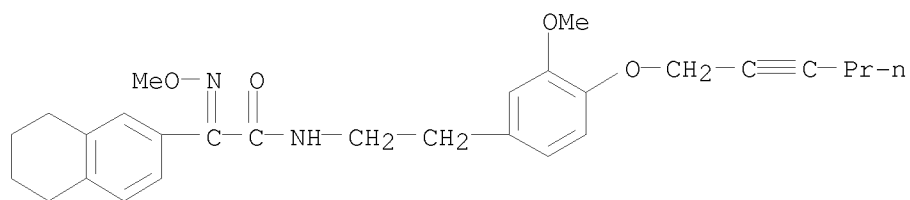


RN 1100623-97-9 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-[(3-cyclopropyl-2-propyn-1-yl)oxy]-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-hydroxy- (CA INDEX NAME)

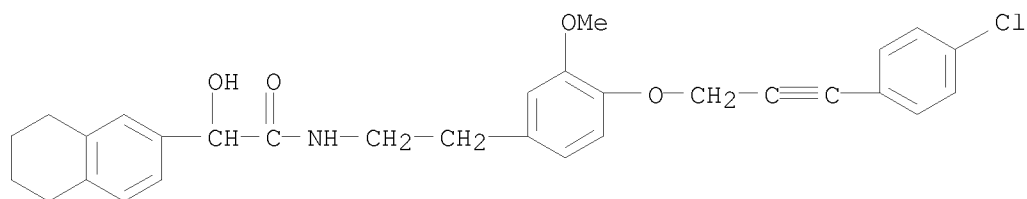


RN 1100626-75-2 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

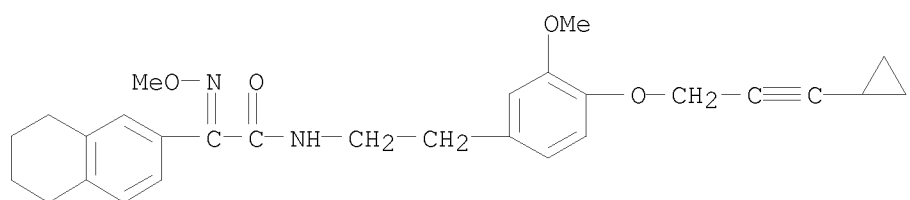
10/513699



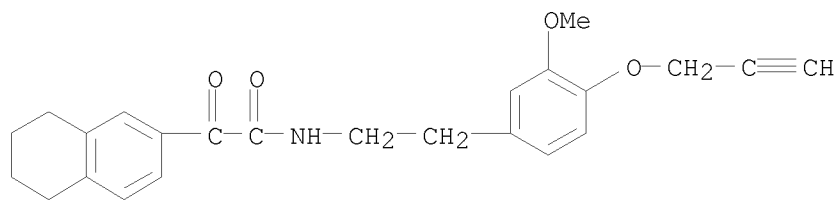
RN 1100628-19-0 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



RN 1100629-73-9 CAPLUS
CN INDEX NAME NOT YET ASSIGNED

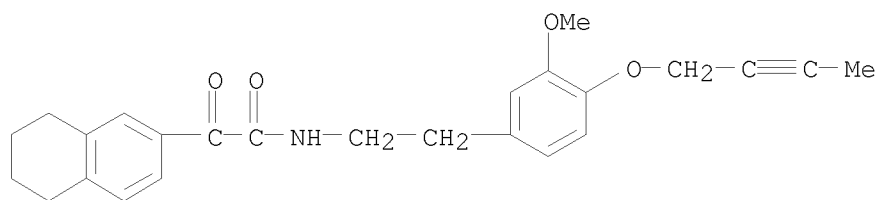


RN 1100630-98-5 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-[2-[3-methoxy-4-(2-propyn-1-yloxy)phenyl]ethyl]-α-oxo- (CA INDEX NAME)

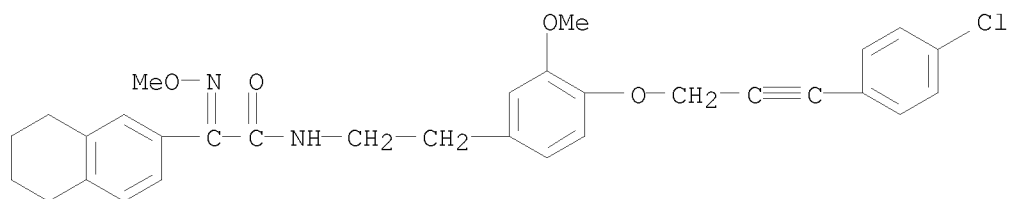


RN 1100633-66-6 CAPLUS
CN 2-Naphthaleneacetamide, N-[2-[4-(2-butyn-1-yloxy)-3-methoxyphenyl]ethyl]-5,6,7,8-tetrahydro-α-oxo- (CA INDEX NAME)

10/513699



RN 1100634-28-3 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



OS.CITING REF COUNT: 11 THERE ARE 11 CAPLUS RECORDS THAT CITE THIS
RECORD (13 CITINGS)
REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS
RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L5 ANSWER 16 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN
 ACCESSION NUMBER: 1965:90683 CAPLUS
 DOCUMENT NUMBER: 62:90683
 ORIGINAL REFERENCE NO.: 62:16162c-e
 TITLE: New tetrahydronaphthalene derivatives
 PATENT ASSIGNEE(S): Holding Ceresia S.A.
 SOURCE: 10 pp.
 DOCUMENT TYPE: Patent
 LANGUAGE: Unavailable
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
FR 1390056		19650219	FR 1964-971752	19640421 <--
BE 658392			BE	
PRIORITY APPLN. INFO.:			CH	19631109

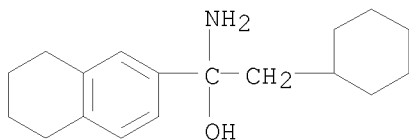
GI For diagram(s), see printed CA Issue.

AB Condensation of ω -bromo-5,6,7,8-tetrahydro-2-acetonaphthone (I) with amines leads to substituted ω -amino-5,6,7,8-tetrahydro-2-acetonaphthones (II). The latter can be reduced by hydrogenation to give III. To a solution of 120 g. I in 700 mL. anhydrous Et₂O was added at 15° 60 g. iso-PrNH₂ in 200 mL. Et₂O. After standing 15-20 h. at 10°, the solution was filtered, evaporated in vacuo, and the residue taken up in Et₂O and treated with HCl gas to give ω -isopropylamino-5,6,7,8-tetrahydro-2-acetonaphthone-HCl (II) (R : iso-Pr) (IV), m. 208-9°. An alc. solution of 50 g. IV was hydrogenated over 2 g. Pd-C (10% Pd) at 2-5 atmospheric, filtered, and evaporated to give 2-isopropylamino-1-(5,6,7,8-tetrahydro-2-naphthyl)ethanol-HCl (III) (R : iso-Pr), m. 160-2°. Similarly prepared were the following compds. (compound, R, and m.p. given): II, sec-Bu, 182-4°; III, sec-Bu, 129-30°; II, tert-Bu, 226-7°; III, tert-Bu, 196-7°; II, PhCH₂CHMe (V), 210-11°; III, PhCH₂CHMe, 142-3°; II cyclohexyl, 231-2°; III, cyclohexyl, 178-9°. Most HCl salts were recrystd. from EtOH-Et₂O, but V was recrystd. from EtOH.

IT 1087756-57-7P
 RL: SPN (Synthetic preparation); PRP (Properties); PREP (Preparation)
 (New tetrahydronaphthalene derivatives)

RN 1087756-57-7 CAPLUS

CN 2-Naphthalenemethanol, α -amino- α -(cyclohexylmethyl)-5,6,7,8-tetrahydro-, hydrochloride (1:1) (CA INDEX NAME)



● HCl

10/513699

<12/04/2007>

Erich Leese

L5 ANSWER 17 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1964:432238 CAPLUS

DOCUMENT NUMBER: 61:32238

ORIGINAL REFERENCE NO.: 61:5580b-d

TITLE: New series of β -adrenergic blocking agents

AUTHOR(S): Ferrari, G.; Casagrande, C.; Canova, M.

CORPORATE SOURCE: Lab. Ric. Simes, Milan, S.p.A.

SOURCE: Bollettino Chimico Farmaceutico (1964),

103(1), 32-6

CODEN: BCFAAI; ISSN: 0006-6648

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

GI For diagram(s), see printed CA Issue.

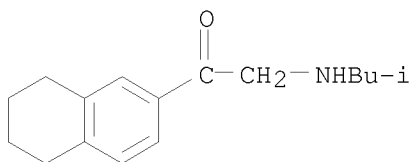
AB The compds. were prepared by catalytic hydrogenation of aminoketones and isolated as hydrochlorides; they were soluble in H₂O and alc., less soluble in Me₂CO, and insol. in ether. The aminoketones used were obtained by reaction, in anhydrous ether, of ω -bromo-5,6,7,8-tetrahydro-2-acetonaphthone (I) (prepared by bromination, in ether, in the presence of Cl₃Al) with an excess of amine containing the iso-Pr, iso-Bu, tert-Bu, cyclohexyl, and 2-phenyl isopropyl radicals. In an example, 160 ml. anhydrous ether containing 25.3 g. I was mixed with 12.7 g. iso-PrNH₂ in 20 ml. anhydrous ether at 15-20°; after continuous stirring 7 hrs., the mixture was filtered, the solvent evaporated in vacuo, the non-reacted amine filtered, and the filtrate acidified with an ether solution of dry HCl to yield 19 g. ω -(isopropylamino)5,6,7,8-tetrahydro-2-acetonaphthone hydrochloride, m. 208-9°. Similarly were prepared the following N-substituted ω -amino5,6,7,8-tetrahydro-2-acetonaphthones (substituent and m.p. given): iso-Bu, 182-4°; tert-Bu, 226-7°; cyclohexyl, 231-2°. and 2-phenylisopropyl, 210-11°. Also prepared were the following (II) (R and m.p. given): iso-Pr, 160-2°; iso-Bu, 129-30°; tert-Bu, 196-8°; cyclohexyl, 178-9°; and 2-phenylisopropyl, 143-4°.

IT 1082682-40-3P

RL: SPN (Synthetic preparation); PRP (Properties); PREP (Preparation)
(New series of β -adrenergic blocking agents)

RN 1082682-40-3 CAPLUS

CN Ethanone, 2-[(2-methylpropyl)amino]-1-(5,6,7,8-tetrahydro-2-naphthalenyl)-
(CA INDEX NAME)



L5 ANSWER 18 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN
 ACCESSION NUMBER: 1964:404138 CAPLUS
 DOCUMENT NUMBER: 61:4138
 ORIGINAL REFERENCE NO.: 61:625d-h,626a-d
 TITLE: Homocyclic compositions
 INVENTOR(S): Howe, R.; Smith, L. H.; Stephenson, J. S.
 PATENT ASSIGNEE(S): Imperial Chemical Industries Ltd.
 SOURCE: 22 pp.
 DOCUMENT TYPE: Patent
 LANGUAGE: Unavailable
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE	
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BE 630210		19631021	BE		<--
FR M3063			FR		
GB 1005026			GB		
NL 290728			NL		
PRIORITY APPLN. INFO.:			GB	19620328	

OTHER SOURCE(S): MARPAT 61:4138

GI For diagram(s), see printed CA Issue.

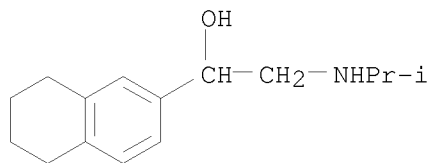
AB Pharmaceutically active compds. of the general formula I where R is a lower alkyl radical, n is 3 or 4, and B may be partially reduced or contain a Me substituent. Thus, NaBH₄ 1 is added with agitation at 0° over 10 min. to 2-(isopropylaminoacetyl)-5,6,7,8-tetrahydronaphthalene-HBr (II) 3 in MeOH 50, the mixture held 3 hrs., the MeOH evaporated at 30°, 0.5N HCl 80 added, the mixture washed with Et₂O 20, 2N NaOH 30 added to the aqueous acidic layer and the aqueous layer extracted with Et₂O 50 parts, dried over anhydrous MgSO₄, and evaporated to give 2-isopropylamino-1-(5,6,7,8-tetrahydro-2-naphthyl)ethanol (III), m. 84-5° (petr. ether, b. 4060°, and AcOEt); HCl salt m. 157°; HBr salt m. 224-6°. 2-[N-(1-Phenylprop-2-yl)amino]-1-(5,6,7,8-tetrahydronaphth-2-yl)ethanol (IV) is produced as a tar in a similar manner starting with the HBr salt of 2-[N-(1-phenylprop-2-yl)aminoacetyl]5,6,7,8-tetrahydronaphthalene. IV oxalate m. 158-9°; IV. HBr m. 227-8°. NaBH₄ 22 is added in 30 min. at 0-15° to 2-(α-bromoacetyl)-5,6,7,8-tetrahydronaphthalene 77, in cyclohexane 200, the mixture kept 1 hr., poured onto ice, and extracted with Et₂O 300 parts, and the extract washed with H₂O, dried over anhydrous MgSO₄, and distilled This oily mixture 10 is heated 16

hrs. with iso-PrNH₂ 20 and EtOH 200, N HCl 100 added, the mixture washed with Et₂O 50, treated with 2N NaOH 75 and extracted with 100 parts Et₂O, the extract washed with H₂O, dried over anhydrous MgSO₄, and evaporated, and the residue treated with oxalic acid in MeOH to give the hemioxalate of 2-isopropylamino-1-(5,6,7,8-tetrahydro-2-naphthyl)ethanol (V), m. 214°. NaBH₄ 1 is added during 10 min. at 0° with stirring to crude (5,6,7,8-tetrahydro-2-naphthyl)glyoxal (VI) in EtNH₂ 3 and MeOH 20, the mixture stirred 2 hrs. and evaporated, the residue treated with 0.5N HCl 100, the solution washed with Et₂O 30, treated with 2N NaOH 35, and extracted with Et₂O 100 parts, the extract washed with water, dried, and evaporated to give 2-ethylamino-1-(5,6,7,8-tetrahydro-2-naphthyl)ethanol (VII), m. 85-6° (AcOEt). VI is obtained by boiling 2-α-bromoacetyl-5,6,7,8-tetrahydronaphthalene S with Me₂SO 70, pouring

onto ice 200, extracting with Et2O 200, washing with saturated NaHCO3 solution 30 parts, then H2O, drying over anhydrous MgSO4, and evaporating 2-Isopropylamino-1-(2-naphthyl)ethanol (VIII) 10 in EtOH 10 is hydrogenated at 125 atmospheric 6 hrs. in the presence of Raney Ni 1, the mixture treated with Et2O 50 and filtered, the filtrate evaporated, the residue treated with 2N HCl 50, the solution washed with Et2O 50, treated with 11N NaOH 20, and extracted with Et2O 50 parts, the extract dried over anhydrous gSO4, evaporated, and recrystd. from petr. ether (b. 60-80°) to give III. VIII 2.3 is hydrogenated at 125°/125 atmospheric 6 hrs. in the presence of 5% Rh-C 0.5, the mixture purified, treated with (CO2H)2 1 in Et2O 50 parts, and filtered to give the oxalate of 1-decahydro-2naphthyl-2-isopropylethanol (IX), m. 122-4° (EtOH-AcOEt, 1:10). Prepared in a similar manner to VII is 2-[N-1-hydroxy-2methylprop-2-yl)amino]-1-(5,6,7,8-tetrahydro-2-naphthyl)ethanol (X), m. 118-19° (AcOEt). Prepared in a similar manner to III are: 1-indan-5-yl-2-isopropylaminoethanol (XI), m. 99°; 2-sec-butylamino-1-indan-5-ylethanol (XII), m. 75-6°; 2-tertbutylamino-1-indan-5-ylethanol (XIII), m. 121-2°; 2-butylaminoindan-5-ylethanol (XIV), m. 94-5°; 2-[2-(3,4-dimethoxyphenyl)ethylamino]-1-indan-5-ylethanol (XV), m. 111-12°; 2tert-butyl-1-(5,6,7,8-tetrahydro-2-naphthyl)ethanol (XVI), m. 85-6° (HCl salt, m. 203.4°). NaBH4 1 is added over 30 min. at 0° to indan-5-ylglyoxal (XVII) 2, EtNH2 1.2, and MeOH 40, the mixture kept 2 hrs. and evaporated, 0.5N HCl 100 added, the mixture washed, with Et2O 30, treated with 2N NaOH 35, and extracted with Et2O 100 parts, and the extract washed with H2O, dried, and evaporated to give 2-ethylamino-1-indan-5-ylethanol (XVIII), m. 110-11° (AcOEt). XVII, m. 240-1° is produced by boiling 5- α -bromoacetylindan 5 and Me2SO 40, keeping 2 days, pouring over ice 200, extracting with Et2O 200, washing with saturated NaHCO3 solution 30 parts, then water, drying over anhydrous MgSO4, evaporating and recrystg. from H2O. A mixture of 2-N-benzyl-N-isopropylamino-1(2-naphthyl)ethanol (XIX) 1, EtOH 16, and concentrated HCl 0.2 is hydrogenated at atm. pressure using Pt oxide 0.3 parts and filtered and the filtrate evaporated, to give 2-isopropylamino-1-(5,6,7,8-tetrahydro-2-naphthyl)ethanol (XX), m. 157° (AcOEt). XIX, m. 154°, is obtained by hydrogenating 2-N-benzyl-N-isopropylaminoacetylnaphthalene using a Pt oxide catalyst. 2-Acetyl-3methyl-5,6,7,8-tetrahydronaphthalene (XXI) 20 is reduced by boiling 5 hrs. with dioxane 150, H2O 10, and SeO2 12.5, cooling, filtering, drying the filtrate, dissolving in iso-PrNH2 7 and EtOH 160, cooling, and adding NaBH4 9 parts in 1 hr. Water 10 is added, the mixture evaporated, the residue washed with Et2O 200 and H2O 50 parts, the ether phase washed with H2O, dried over anhydrous MgSO4, and evaporated to give 2-isopropylamino-1-(3-methyl5,6,7,8-tetrahydro-2-naphthyl)ethanol (XXII), m. 108° (Ac-OEt). XXI, b8 153-7° is obtained by adding a mixture of 2-methyl-5,6,7,8-tetrahydronaphthalene 15, AcCl 9, and CS2 60 at 0° to a suspension of AlCl3 15 in CS2 125 parts. After 16 hrs. an ice-H2O mixture 200 is added, the CS2 evaporated, the residue extracted with Et2O 200, the extract washed with H2O, dried over anhydrous MgSO4, and evaporated, and the residual liquid fractionally distilled Tablets for oral administration are obtained from III.HCl, IV.HCl, VII.HCl, X.HCl, or XIII.HCl.

10/513699

IT 1071607-66-3P
RL: SPN (Synthetic preparation); PRP (Properties); PREP (Preparation)
(Homocyclic compositions)
RN 1071607-66-3 CAPLUS
CN 2-Naphthalenemethanol, 5,6,7,8-tetrahydro- α -[[1-methylethyl)amino]methyl]-, hydrobromide (1:1) (CA INDEX NAME)



● HBr

L5 ANSWER 19 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1962:423372 CAPLUS

DOCUMENT NUMBER: 57:23372

ORIGINAL REFERENCE NO.: 57:4722c-i, 4723a-c

TITLE: Chemistry of p-quinols. I. V. Stereochemistry of the Tetralin p-quinols and the estra-p-quin-10-ols

AUTHOR(S): Hecker, Erich; trell, Rudolf Lat; Meyer, Elisabeth

CORPORATE SOURCE: Biochem., Max-Planck-Inst., Munich, Germany

SOURCE: Chemische Berichte (1962), 95, 985-95

CODEN: CHBEAM; ISSN: 0009-2940

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

AB Racemic Tetralin p-quinol (I) was resolved into the optical antipodes via the strychnine (II) salts of the acid 3-nitrophthalates. The configuration and conformation of the antipodes is discussed and related to the configuration and conformation of the estra-p-quinols on the basis of the optical shift rule and the rotational dispersion. From this the absolute configuration of the antipodes of I is deduced. Tetralin nitrated by the method of Schroeter (CA 16, 1763) gave a yellow oily mixture of 5- and 6-nitro derivs., b13 156-65°; a 90-g. portion in 400 cc. hot EtOH treated with 5.5 g. NH₄Cl in 80 cc. H₂O and then with stirring above 70° during 10 min. with 70 g. Zn dust, stirred 10 min., filtered at 65-70°, the residue washed with 80% EtOH, the combined filtrates poured into 2.5 l. H₂O and extracted with 600 cc. Et₂O, the extract shaken immediately with 250 cc. 10% H₂SO₄ to precipitate the sulfate of the 6HONH derivative

of Tetralin, the aqueous phase again extracted with 600 cc. Et₂O, and the extract

shaken with H₂SO₄ and filtered, the Et₂O phase separated, the filter residues suspended in the combined aqueous phases and extracted 10 hrs. with C₆H₆, and

the

extract evaporated gave 22 g. dark oil which diluted with 50 cc. cyclohexane deposited 5-6 g. pure I, m. 125-6° (Me₂CO and EtOAc); the original Et₂O phase evaporated, the residue (55 g.) distilled, and the distillate (40

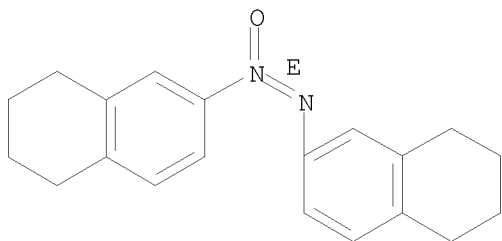
g.),

b13 150-65°, refrigerated gave 5-nitro derivative of Tetralin, m. 34-5° (MeOH); the distillation residue (15 g.) gave some 6,6'-azoxy derivative of Tetralin, m. 99-100°. I (1.3 g.) and 4.3 g. 3-nitrophthalic anhydride heated 7 hrs. at 40-5° in 14 cc. CH₂Cl₂ and 14 cc. C₅H₅N, kept overnight, heated 1 hr. at 40°, poured into iced H₂O, acidified with N HCl, and extracted with 3:1 CHCl₃-CH₂Cl₂, and the extract worked up gave 1.73 g. 3-nitro-2-phthalate (III) of I, m. 172° (MeOH); the mother liquor evaporated, and residue (1.6 g.) (from several runs) dissolved in 2.5 cc. warm MeOCH₂CH₂OH and allowed to stand 2 days deposited 820 mg. 3-nitro-1-phthalate (IV) of I, m. 192°. The R_f values were determined with 15:5:6 nonane-C₆H₆-AcOH for the following compds. (m.p. and R_f value given): Et 3-nitro-2 phthalate, 154°, 0.24; Et 3-nitro-1-phthalate, 107°, 0.18; Bu 8-nitro-2-phthalate, 143°, 0.40; Bu 3-nitro-1-phthalate, 90°, 0.34; III, 172°, 0.18; IV, 192°, 0.13. III (7 mg.) treated 10 hrs. at 45° with a 5-fold excess of K₂CO₃ gave only. unchanged III. II (7 mg.) refluxed 7 hrs. with 2.5 equivs. KOH in MeOH gave only phenolic material and unchanged III; the same result was obtained by refluxing 4 hrs. with 1% H₂SO₄-MeOH or by keeping several days in MeOH with 1.5 equivs. p-MeC₆H₄SO₃H. III (1.067 g.) in 100 cc. CHCl₃ treated with CH₂N₂-Et₂O and the product chromatographed on Al₂O₃ gave 95% 1-Me 2-(Tetralin pquinol) 3-nitrophthalate (V), m. 114-15.5° (1:1

C₆H₆cyclohexane). Crude V (3 millimoles) in 25 cc. MeOH treated with 60 cc. aqueous K₂CO₃, kept 5 days at room temperature, treated in the dark with H₂O and CHCl₃, and the residue from the organic layer chromatographed on Al₂O₃ yielded 100 mg. yellow, partially crystalline mixture of polymethylene and diMe 3-nitrophthalate (VI), 283 g. crystalline VI, 250 mg. V, 40 mg. yellow oily mixture of III and I, and 140 mg. I. III (7.14 g.) in 250 cc. Me₂CO treated with 6.68 g. II in the min. amount of CHCl₃, concentrated in vacuo at 40° to about 100 cc., diluted with 150 cc. Me₂CO, again concentrated to 100 cc., diluted with 40 cc. H₂O of 60°, concentrated to 120 cc., and refrigerated overnight gave 6.2 g. II salt; the salt (6.2 g.) in 20 cc. CHCl₃ diluted with 200 cc. Me₂CO and concentrated to 100 cc. at 40°, this treatment repeated, and the concentrate diluted with 40 cc. warm H₂O and refrigerated overnight gave 5.4 g. salt, [α]_D²⁰ -12° (4:1 EtOH-CHCl₃), which recrystd. gave 4.3 g. salt, [α]_D²⁵ -11° (4:1 EtOH-CHCl₃). II salt (4.5 g.), [α]_D²⁴ -12°, in 100 cc. CHCl₃ shaken with five to seven 30cc. portions 2N HCl, concentrated to 4 cc., and filtered on the next day gave (+)-III, [α]_D²⁵ 32° (c 1.25, EtOH), [α]_D²⁶ 62° (c 1.25, dioxane), m. 173-4° (MeOH). The original mother liquor evaporated, the residual II salt (7.5 g.), [α]_D²² 12° (EtOH-CHCl₃), dissolved in 20 cc. H₂O, and slowly evaporated gave 2.1 g. crystals, [α]_D²⁴ 8° (EtOH-CHCl₃); the remaining mother liquor gave 5.6 g. II salt, [α]_D²⁶ 14° (EtOH-CHCl₃); a 5-g. portion decomposed with 2N HCl yielded 1.81 g. (-)-III, m. 170-1° (CHCl₃), [α]_D²⁶ -30° (c 0.93, EtOH), [α]_D²⁵ - 62° (c 1.69, dioxane). (+)-III (1.07 g.) treated with CH₂N₂-Et₂O, the resulting Me ester saponified with K₂CO₃ in MeOH, and the reaction product chromatographed gave 110 mg. (+)-V, m. 119-20°, [α]_D²⁵ 114°, [α]_D²⁶ 111° (c 1, dioxane). (-)-III (1.8 g.) gave similarly 92 mg. (-)-V, m. 118-19° (EtOAc), [α]_D²⁰ - 12° (c 1, dioxane).

IT 1087735-80-5P
 RL: SPN (Synthetic preparation); PRP (Properties); PREP (Preparation)
 (Chemistry of p-quinols. I. V. Stereochemistry of the Tetralin
 p-quinols and the estra-p-quin-10-ols)
 RN 1087735-80-5 CAPLUS
 CN INDEX NAME NOT YET ASSIGNED

Double bond geometry as shown.



OS.CITING REF COUNT: 1 THERE ARE 1 CAPLUS RECORDS THAT CITE THIS RECORD
 (1 CITINGS)

L5 ANSWER 20 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1960:97472 CAPLUS

DOCUMENT NUMBER: 54:97472

ORIGINAL REFERENCE NO.: 54:18451c-h

TITLE: Derivatives of naphthalenes and fatty acids:
heptanoylnaphthalenes and -naphthols

AUTHOR(S): Jorand, J.

SOURCE: Oleagineux (1960), 15, 183-8
CODEN: OLEAAF; ISSN: 0030-2082

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

- AB The following derivs. of Me(CH₂)₅CO₂H, b₂ 134°, were prepared:
 Me(CH₂)₅COC1 (I), b₂₅₋₂₇ 83-6°; Me(CH₂)₅CONH₂, m. 96-6.5°;
 Me(CH₂)₅CN (II), b_{2885°}, n_{20D} 1.4144; α-C₁₀H₇CO(CH₂)₅Me (III)
 (from II and α-C₁₀H₇MgBr), b_{0.01-0.02} 162-8°, n_{20D} 1.5715,
 n_{40D} 1.5640; III 2,4-dinitrophenylhydrazone m. 128.5;
 β-C₁₀H₇CO(CH₂)₅Me (IV) (by Friedel-Crafts), m. 60°, b_{0.02}
 155-8°; IV 2,4-dinitrophenylhydrazone m. 191.5-2°; IV
 p-nitrophenylhydrazone m. 174-5°; β-heptanoyl derivative (V) of
 Tetralin, b_{0.02} 152-4°, n_{20D} 1.5320, n_{40D} 1.5241; V
 2,4-dinitrophenylhydrazone m. 155.5°. The absorption spectra of
 the different dinitrophenylhydrazones were given. Also prepared were
 α-C₁₀H₇CH(OH)(CH₂)₅Me, n_{20D} 1.5692, n_{40D} 1.5609;
 β-C₁₀H₇CH(OH)(CH₂)₅Me, m. 35-7°; acetate m. 60°, n_{40D}
 1.5537; β-C₁₀H₁₁CH(OH)(CH₂)₅Me, n_{20D} 1.5241, n_{40D} 1.5172;
 α-C₁₀H₇CH(NH₂)(CH₂)₅Me, b_{0.02} 152-5°, n_{20D} 1.5702;
 hydrochloride m. 189.5-90.5°; picrate m. 201°;
 phenylthiourea derivative m. 141.5-2.0°; β-C₁₀H₇CH(NH₂)(CH₂)₅Me,
 b_{0.01} 168-70°, n_{20D} 1.5700; hydrochloride m. 199-200°;
 picrate m. 208-8.5°; phenylthiourea derivative m. 114-15°;
 β-C₁₀H₁₁CH(NH₂)(CH₂)₅Me, b_{0.02} 136-8°, n_{20D} 1.5720;
 hydrochloride m. 205.5-6.5°; styphnate m. 201-2°. I (75 g.)
 added at 10° to 72 g. α-C₁₀H₇OH and 70 g. ZnCl₂ in 300 cc.
 C₆H₅NO₂, the mixture kept 48 hrs. at room temperature, hydrolyzed with HCl-ice,
 the C₆H₅NO₂ layer decanted, distilled in vacuo, the residue extracted with 10%
 KOH solution, the fraction insol. in alkali washed with H₂O, and the product
 crystallized from 95% EtOH gave 1-HOC₁₀H₆CO(CH₂)₅Me-2 (VI), m. 52°
 (MeOH); 2,4-dinitrophenylhydrazone m. 216°. Also prepared were
 2-HOC₁₀H₆CO(CH₂)₅Me-1 (VII); C₁₀H₇OCO(CH₂)₅Me-1 (from α-C₁₀H₇OH and
 I in C₆H₆pyridine), b_{0.02} 143-6°; C₁₀H₇OCO(CH₂)₅Me-2 (VIII), b_{0.02}
 157-65°, m. 38°. VII (by Fries rearrangement of VIII) b_{0.02}
 150-5°, n_{20D} 1.5922; 2,4-dinitrophenylhydrazone could not be
 obtained. AlCl₃ (75 g.) added to 75 g. I and 79 g. 1-C₁₀H₇OMe in 400 cc.
 PhNO₂ with vigorous stirring at 0°, the complex hydrolyzed,
 neutralized, PhNO₂ distilled, and the raw product distilled in vacuo gave
 1-MeOC₁₀H₆CO(CH₂)₅Me-4 (IX), b_{0.02-0.05} 200-5°, m. 43°
 (MeOH); 2,4-dinitrophenylhydrazone m. 125.5. IX demethylated according to
 Buu-Hoi (CA 44, 4444f) at 200° with pyridine-HCl gave
 1-HOC₁₀H₆CO(CH₂)₅Me-4, m. 116°; 2,4-dinitrophenylhydrazone m.
 202°. 2-MeOC₁₀H₆CO(CH₂)₅Me-6 m. 70°, b_{0.02-0.05}
 170-80°; 2,4-dinitrophenylhydrazone m. 189°.
 2-HOC₁₀H₆CO(CH₂)₅Me-6 m. 135-5.5°; 2,4-dinitrophenylhydrazone m.
 251°.
- IT 860221-29-0P, Styphnic acid, compound with
 α-hexyl-5,6,7,8-tetrahydro-2-naphthalenemethylamine
 RL: PREP (Preparation)
 (preparation of)

10/513699

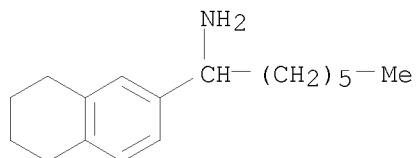
RN 860221-29-0 CAPLUS

CN 2-Naphthalenemethanamine, α -hexyl-5,6,7,8-tetrahydro-, compd. with
2,4,6-trinitro-1,3-benzenediol (1:1) (CA INDEX NAME)

CM 1

CRN 101744-83-6

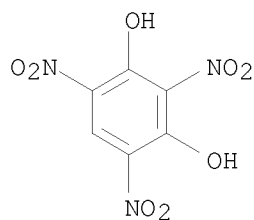
CMF C17 H27 N



CM 2

CRN 82-71-3

CMF C6 H3 N3 O8



OS.CITING REF COUNT: 1

THERE ARE 1 CAPLUS RECORDS THAT CITE THIS RECORD
(1 CITINGS)

L5 ANSWER 21 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1960:50301 CAPLUS

DOCUMENT NUMBER: 54:50301

ORIGINAL REFERENCE NO.: 54:9854c-g

TITLE: Possible antituberculous compounds. VII. Preparation of 5,6,7,8-tetrahydro-1(and 2)-naphthylamidines

AUTHOR(S): Misra, Vinay S.; Husain, Md. Imtiaz

CORPORATE SOURCE: Univ. Lucknow

SOURCE: Journal of the Indian Chemical Society (1959), (36), 803-6

CODEN: JICSAH; ISSN: 0019-4522

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

AB cf. C.A. 50, 15476g. The observation of Oxley and Peak (cf. Charlton, et al., C.A. 46, 2005a) that N-(1-naphthyl)benzamidine (I) had much greater antitubercular activity than α -naphthylamine (II) led M. and H. to prepare the title compds. whose surface area and lipoid solubility were greater than those of I. II (10 g.) in 112 g. boiling amyl alc. was added in a continuous stream to 8 g. Na, the mixture heated on a steam bath until all the Na had disappeared, the product poured into water, the upper layer separated, the bases converted into HCl salts, dissolved in hot water, treated with excess NaOH, and the organic layer washed and distilled to give 4.5 g. 5,6,7,8-tetrahydro-1-naphthylamine (III), b2-3 135°; Ac derivative m. 158° (water). β -Naphthylamine (10 g.) was similarly reduced to 4 g. 5,6,7,8-tetrahydro-2-naphthylamine (IV), b2 160°. A stirred solution of 4.0 g. III in 150 cc. dry Et2O was treated dropwise with 5 g. PhSO3H in MeOH, the product filtered off, and crystallized from hot water to give 7.5 g. 5,6,7,8-tetrahydro-1-naphthyl-ammonium benzenesulfonate (V), m. above 275°, dark brown amorphous powder. IV (4.0 g.) and 5 g. PhSO3H gave 3.0 g. 5,6,7,8-tetrahydro-2-naphthylammonium benzenesulfonate, m. 194°, light brown amorphous powder. The sulfonates were fused at 230-5° with the required nitriles to give the desired amidinium benzenesulfonates; subsequent treatment with base gave the free amidines (aryl group, % yield, m.p., color, m.p. of benzenesulfonate, its % yield, and color given): Ph, 86, 95-6°, brown, 235-6°, 23, yellow; o-tolyl, 79, 246°, white, 260-1°, 22, brown; m-tolyl, 6, 180°, dark brown, - (would not crystallize), -, -; p-tolyl, 6, above 280°, brown, - (would not crystallize), -, -; α -naphthyl, 77, 210°, black, 270°, 26, black. The corresponding 2-amidines (same data given): o-tolyl, 90, 115°, brown, above 285°, 10%, brown; m-tolyl, 46, above 290°, light green, - (would not crystallize), -, -; p-tolyl, 80, 150°, dark brown, above 270°, 28, brown; α -naphthyl, 64, above 280°, black, above 270°, 48, black; Ph, -, 285°, 12, white.

IT 853648-64-3P, 2-Naphthylamine, 5,6,7,8-tetrahydro-, compound with benzenesulfonic acid

RL: PREP (Preparation)
(preparation of)

RN 853648-64-3 CAPLUS

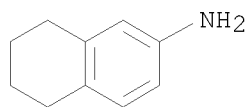
CN 2-Naphthalenamine, 5,6,7,8-tetrahydro-, benzenesulfonate (1:1) (CA INDEX NAME)

CM 1

CRN 2217-43-8

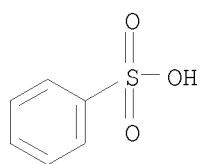
CMF C10 H13 N

10/513699



CM 2

CRN 98-11-3
CMF C6 H6 O3 S



<12/04/2007>

Erich Leese

10/513699

L5 ANSWER 22 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1960:45099 CAPLUS

DOCUMENT NUMBER: 54:45099

ORIGINAL REFERENCE NO.: 54:8954e-h

TITLE: Liberation of bradykinin from plasma by treatment with
peptone or by boiling with hydrochloric acid

AUTHOR(S): Rocha e Silva, M.; Holzhacker, Elisabeth L.

CORPORATE SOURCE: Inst. biol., Sao Paulo, Brazil

SOURCE: Archives Internationales de Pharmacodynamie et de
Therapie (1959), 122, 168-79
CODEN: AIPTAK; ISSN: 0003-9780

DOCUMENT TYPE: Journal

LANGUAGE: English

AB Different fractions of peptone (proteose-peptone, Difco) were obtained by chromatography on the ion exchange resin Amberlite IRC-50. A fraction eluted at pH 7.8 (1.5 g. from about 30 g. of peptone) showed the greatest activity in releasing bradykinin (I) when incubated with heparinized rat plasma. It was also the most active fraction as a histamine releaser. Plasma from rats, when heated 1-10 min. with 0.1N HCl, developed full I activity upon neutralization and incubation at room temperature or at 37°. This spontaneous release of I was enzymic in nature and was inhibited by soy-bean trypsin inhibitor. The material released by HCl or by the peptone fraction was identified as I by parallel assays on the guinea pig ileum, on the uterus of the rat (the most sensitive method), or by effects of the blood pressure of rabbits. Since the release of I from denatured plasma by trypsin and by snake venom is parallel to the resp. esterase activities of these agents against benzoyl-L-arginine methyl ester, the release of I might provide a very sensitive indication of activation of an enzyme in plasma displaying a similar activity. 18 references.

IT 856631-76-0P, Benamidine, N-[5,6,7,8-tetrahydro-2-naphthyl]-, compds. with benzenesulfonic acid

RL: PREP (Preparation)

(preparation of)

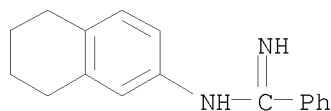
RN 856631-76-0 CAPLUS

CN Benzenecarboximidamide, N-(5,6,7,8-tetrahydro-2-naphthalenyl)-, benzenesulfonate (1:1) (CA INDEX NAME)

CM 1

CRN 109249-02-7

CMF C17 H18 N2

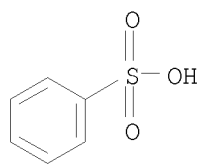


CM 2

CRN 98-11-3

CMF C6 H6 O3 S

10/513699



<12/04/2007>

Erich Leese

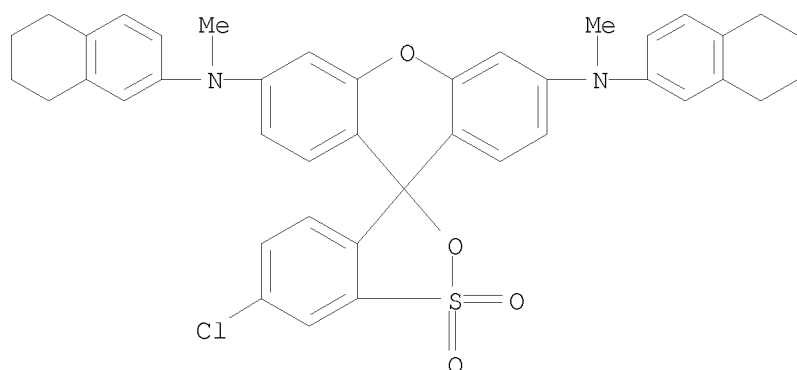
L5 ANSWER 23 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1956:75767 CAPLUS
 DOCUMENT NUMBER: 50:75767
 ORIGINAL REFERENCE NO.: 50:14237f-i,14238a-b
 TITLE: Triphenylmethane dyes
 PATENT ASSIGNEE(S): Farberke Hoechst AG vorm. Meister Lucius & Bruning
 DOCUMENT TYPE: Patent
 LANGUAGE: Unavailable
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
	-----	----	-----	-----	-----
	GB 744972		19560215	GB 1951-17267	19510720 <--
AB	3,6-Diamino-9-(2-sulfophenyl)xanthene dyes, whose amino groups can be substituted by aliphatic or aromatic radicals or whose N may be a member of a heterocyclic ring, are prepared To improve their solubility, the dyes can be further sulfonated. They dye wool and silk in red to blue tints of very good fastness to light and alkalies. 3,6-Dihydroxy-9-(2-sulfophenyl)xanthene (I) 10 is heated with POC13, the excess POC13 is removed by distillation, and the resulting yellow 3,6-dichloro-9-(2-sulfophenyl)xanthene (II) is treated with PhNH2 30 parts several hrs. at 100°. The excess PhNH2 is extracted from the melt with HCl and the dye isolated. After drying it is sulfonated with 95% H2SO4 until it dissolves in dilute Na2CO3, and isolated by pouring into H2O, filtering, dissolving in dilute Na2CO3, and salting out. It dyes a clear violet. The following shades are obtained by replacing PhNH2: red-violet with o-toluidine (III); bright-red with 2,6-Me2C6H3NH2, 2-MeC6H4NH2, 2-MeC6H4NHMe, 2,5-Me2C6H4NHMe, (HOCH2CH2)2NH, or piperidine; reddish blue with 4-EtOC6H4NH2 and a still bluer shade with 2-naphthylamine (IV); violet with 2-NaO3SC6H4OC6H4NH2-4; reddish violet with 2-HO3SC6H4NH2 in glycol; red-violet with 3-ClC6H4NH2; clear violet with 2-MeOC6H4NH2, 2-EtOC6H4NH2, 2,4-H2N(Cl)C6H4OC6H5(V), or 1-naphthylamine (VI); and navy blue by replacing PhNH2 with 2,3-HO(HOOC)C6H3NH2 or 2,3,5-HO(HOOC)(HO3S)C6H2NH2 in glycol but without after-sulfonation. I 10 is heated to 130° with 2,4-Me2C6H3NH2 50 in the presence of POC13 10 parts, heated till the color no longer deepens, and the dye is isolated by dissolving out the amine. It dyes wool violet. Similar dyes are also obtained by using 3,6-dichloro-9-(2,4-disulfophenyl)xanthene (VII) instead of II. VII gives with III, after sulfonation, a soluble clear red-violet dye, with m-toluidine a navy-blue dye, and similar violet shades with VI, various xylidines, chloro- and bromoanilines. A greenish dye is made from VII with IV or 2,3-HO(HOOC)C6H3NH2, blue to violet shades with V, 3,4-HOOC(HO)C6H3NH2 or 2,5,4-Me2(HO3S)C6H2NH2, red-violet to blue-violet shades with 2,4,5-Me2(HO3S)C6H2NH2 or 2,4,6-Me2(HO3S)C6H2NH2, a gray-blue with 4-H2NC6H4NHC6H5, and a red with 4(or 5)-chloro-2-aminobenzotrifluoride. Similar dyes are also obtained by the same process from II substituted in the benzene ring by a p-ethoxy, a p-chloro-, a 4-hydroxy, and a 3-carboxylic acid group.				
IT	853780-16-2P, Benzenesulfonic acid, 5-chloro-2-[9-hydroxy-3,6-bis[methyl(5,6,7,8-tetrahydro-2-naphthyl)amino]-9-xanthenyl]-, γ-sultone RL: PREP (Preparation) (preparation of)				
RN	853780-16-2 CAPLUS				
CN	Spiro[3H-2,1-benzoxathiole-3,9'-[9H]xanthene]-3',6'-diamine, 6-chloro-N3',N6'-dimethyl-N3',N6'-bis(5,6,7,8-tetrahydro-2-naphthalenyl)-,				

10/513699

1,1-dioxide (CA INDEX NAME)



<12/04/2007>

Erich Leese

L5 ANSWER 24 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1956:40401 CAPLUS

DOCUMENT NUMBER: 50:40401

ORIGINAL REFERENCE NO.: 50:7803c-f

TITLE: Chloromethylation of tetralin

AUTHOR(S): Vanags, G.; Gudriniece, E.

SOURCE: Latvijas PSR Zinatnu Akademijas Vestis (1955

), (No. 5 (Whole No. 94)), 119-24

CODEN: LZAVAL; ISSN: 0132-6422

DOCUMENT TYPE: Journal

LANGUAGE: Russian

AB Tetralin (66 mg.), 28 g. (CH₂O)_n, 65 ml. glacial AcOH, 33 g. crystalline H₃PO₄, and 91 ml. concentration HCl at 85-90° stirred 4 hrs. gave 66%

1,2,3,4-tetrahydro-6-chloromethylnaphthalene (I). With excess II, 10% 5,8-bis(chloromethyl)-1,2,3,4-tetrahydronaphthalene was obtained in addition to I. The 6-piperidinomethyl analog (II of I) was prepared by treating I in Et₂O with piperidine at room temperature II decomposed on distillation

Bubbling dry HCl

through II in Et₂O gave II.HCl, very hygroscopic. II picrate, m. 150°. 1-(1,2,3,4-Tetrahydro-6-naphthylmethyl)pyridinium chloride, m. 115°, was prepared (88.5% yield) from 7.2 g. I, 20 ml. absolute Et₂O, and dry pyridine. H₂NC(SR):NH.HCl (R = 1,2,3,4-tetrahydro-6-naphthylmethyl), m. 212°, was prepared (96% yield) by heating 7.2 g. I with 6 g. thiourea. RCO₂H was prepared (42% yield) refluxing crude I with KCN in H₂O, and hydrolyzing the nitrile with aqueous NaOH; the hydrolysis was aided, and formation of resinous products was minimized by adding small amts. of 3% H₂O₂ at intervals. RCONHPh, m. 112°, was obtained by method similar to that described (C.A. 50, 271f).

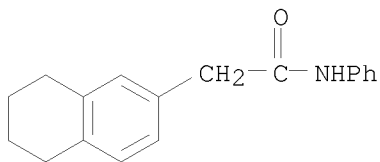
IT 860396-36-7P, 2-Naphthaleneacetanilide, 5,6,7,8-tetrahydro-

RL: PREP (Preparation)

(preparation of)

RN 860396-36-7 CAPLUS

CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-phenyl- (CA INDEX NAME)



L5 ANSWER 25 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1956:10063 CAPLUS
 DOCUMENT NUMBER: 50:10063
 ORIGINAL REFERENCE NO.: 50:2116a-d
 TITLE: 1-Aryl derivatives of
 2-nitro-1-(5,6,7,8-tetrahydro-2-naphthyl)alkanes and
 their use as insecticides
 INVENTOR(S): Johnson, Arnold N.
 PATENT ASSIGNEE(S): Commercial Solvents Corp.
 DOCUMENT TYPE: Patent
 LANGUAGE: Unavailable
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 2719810		19551004	US	<--

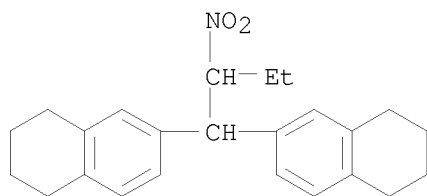
AB Compds. of the type XCH(Ar)CH(NO₂)R, in which X is
 5,6,7,8-tetrahydro-2-naphthyl, R is Me or Et, and Ar is a substituted Ph
 group, were prepared, emulsified in xylene, and tested for insecticidal
 properties by spraying on flies and on beans which were then fed to the
 Mexican bean beetle, the Southern army worm, and the pea aphid. In a
 typical preparation, 41 g. 2-nitro-1-p-tolyl-1-propanol was added to a mixture
 of 102 g. 1,2,3,4-tetrahydronaphthalene and 100 ml. concentrated H₂SO₄ in 20 min.
 at 20-5°. The mixture was agitated for 1 hr. The top layer was
 steam distilled to yield 49.5 g. 2-nitro-1-p-tolyl-1-(5,6,7,8-tetrahydro-2-
 naphthyl)propane, a thick oily product. Six g. of this product recrystd.
 from petroleum hexane, then from EtOH to give 0.7 g. of a white solid, m.
 114-15°. In a similar manner, the following compds. were prepared
 and tested: 2-nitro-1-(p-chlorophenyl)-1-(5,6,7,8-tetrahydro-2-
 naphthyl)butane, 2-nitro-1-(p-chlorophenyl)-1-(5,6,7,8-tetrahydro-2-
 naphthyl)propane, 2-nitro-1-p-tolyl-1-(5,6,7,8-tetrahydro-2-
 naphthyl)butane, 2-nitro-1-(p-methoxyphenyl)-1-(5,6,7,8-tetrahydro-2-
 naphthyl)butane, 2-nitro-1-(3,4-methylenedioxyphenyl)-1-(5,6,7,8-
 tetrahydro-2-naphthyl)butane, 2-nitro-1,1-bis(5,6,7,8-tetrahydro-2-
 naphthyl)butane, 2-nitro-1-(p-ethylphenyl)-1-(5,6,7,8-tetrahydro-2-
 naphthyl)butane, 2-nitro-1-(p-isopropylphenyl)-1-(5,6,7,8-tetrahydro-2-
 naphthyl)butane, 2-nitro-1-(diethylphenyl)-1-(5,6,7,8-tetrahydro-2-
 naphthyl)butane, and 2-nitro-1-xylyl-1-(5,6,7,8-tetrahydro-2-
 naphthyl)butane.

IT 854459-68-0, Butane, 2-nitro-1,1-bis(5,6,7,8-tetrahydro-2-
 naphthyl)- 855952-07-7, Anisole,
 p-[2-nitro-1-(5,6,7,8-tetrahydro-2-naphthyl)butyl]- 858199-39-0
 , Naphthalene, 1,2,3,4-tetrahydro-6-[p-isopropyl- α -(1-
 nitropropyl)benzyl]- 858199-45-8, Naphthalene,
 1,2,3,4-tetrahydro-6-[p-methyl- α -(1-nitropropyl)benzyl]-
 858458-47-6, Naphthalene, 6-[p-ethyl- α -(1-
 nitropropyl)benzyl]-1,2,3,4-tetrahydro- 858459-74-2,
 Naphthalene, 1,2,3,4-tetrahydro-6-[α -(1-nitropropyl)piperonyl]-
 858459-78-6, Naphthalene, 1,2,3,4-tetrahydro-6-(p-methyl- α -1-
 nitroethylbenzyl)- 860366-28-5, Naphthalene,
 6-(p-chloro- α -1-nitroethylbenzyl)-1,2,3,4-tetrahydro-
 860395-78-4, Naphthalene, 6-[p-chloro- α -(1-
 nitropropyl)benzyl]-1,2,3,4-tetrahydro-
 (insecticide)

RN 854459-68-0 CAPLUS

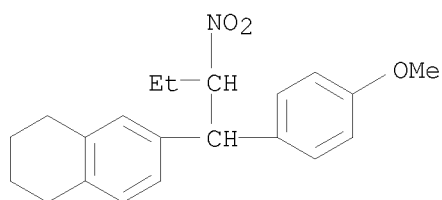
10/513699

CN Butane, 2-nitro-1,1-bis(5,6,7,8-tetrahydro-2-naphthyl)- (5CI) (CA INDEX NAME)



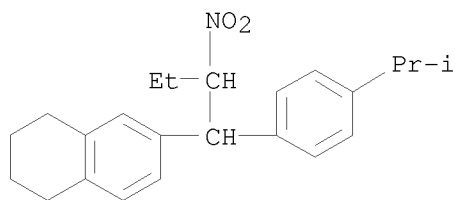
RN 855952-07-7 CAPLUS

CN Naphthalene, 1,2,3,4-tetrahydro-6-[1-(4-methoxyphenyl)-2-nitrobutyl]- (CA INDEX NAME)



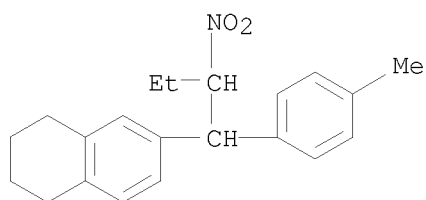
RN 858199-39-0 CAPLUS

CN Naphthalene, 1,2,3,4-tetrahydro-6-[1-[4-(1-methylethyl)phenyl]-2-nitrobutyl]- (CA INDEX NAME)



RN 858199-45-8 CAPLUS

CN Naphthalene, 1,2,3,4-tetrahydro-6-[1-(4-methylphenyl)-2-nitrobutyl]- (CA INDEX NAME)

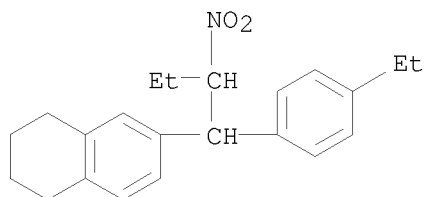


RN 858458-47-6 CAPLUS

CN Naphthalene, 6-[1-(4-ethylphenyl)-2-nitrobutyl]-1,2,3,4-tetrahydro- (CA INDEX NAME)

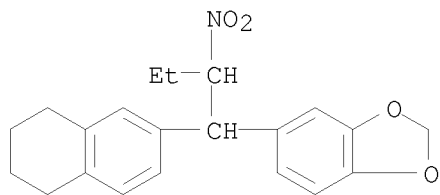
10/513699

INDEX NAME)



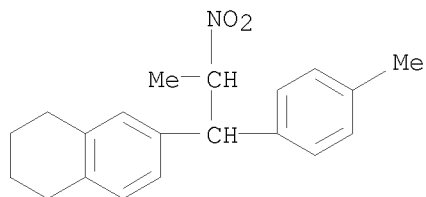
RN 858459-74-2 CAPLUS

CN 1,3-Benzodioxole, 5-[2-nitro-1-(5,6,7,8-tetrahydro-2-naphthalenyl)butyl]-
(CA INDEX NAME)



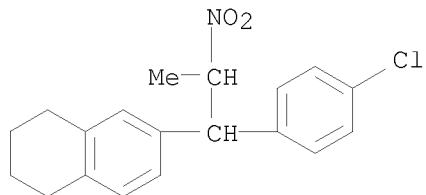
RN 858459-78-6 CAPLUS

CN Naphthalene, 1,2,3,4-tetrahydro-6-[1-(4-methylphenyl)-2-nitropropyl]- (CA
INDEX NAME)



RN 860366-28-5 CAPLUS

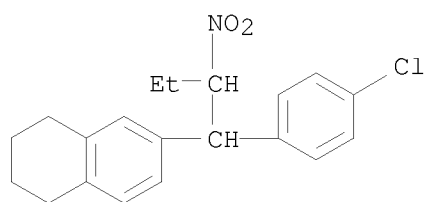
CN Naphthalene, 6-[1-(4-chlorophenyl)-2-nitropropyl]-1,2,3,4-tetrahydro- (CA
INDEX NAME)



RN 860395-78-4 CAPLUS

CN Naphthalene, 6-[1-(4-chlorophenyl)-2-nitrobutyl]-1,2,3,4-tetrahydro- (CA
INDEX NAME)

10/513699



<12/04/2007>

Erich Leese

L5 ANSWER 26 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1955:56624 CAPLUS

DOCUMENT NUMBER: 49:56624

ORIGINAL REFERENCE NO.: 49:10906e-i,10907a-d

TITLE: Use of 1,3-dichloro-2-butene for the synthesis of some ketone derivatives of bicyclo[1.3.3]nonene and of hexahydronaphthalene

AUTHOR(S): Julia, Sylvestre A.

CORPORATE SOURCE: Ecole polytech., Paris Ve

SOURCE: Bulletin de la Societe Chimique de France (1954) 780-9

CODEN: BSCFAS; ISSN: 0037-8968

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

OTHER SOURCE(S): CASREACT 49:56624

AB The condensation of cyclohexanone and its derivs. with 1,3-dichloro-2-butene (I) and the subsequent cyclization of these products was studied with particular reference to steric effects and orientation. 2-Methylcyclohexanone (12 g.) and 12.5 g. I in 50 cc. C₆H₆ agitated and cooled with ice, 2N Na amylate added slowly, the mixture kept at room temperature

for 1 hr., and then refluxed for 3 hrs. gave

2-methyl-2-(γ -chlorocrotyl)cyclohexanone (II), b₁₄ 130°, n_{17D}

1.4941; semicarbazone, m. 153-5°; 2,4-dinitrophenylhydrazone, m.

133-5°. 1,4-Dimethylbicyclo[3.3.1]non-3-en-9-one, b₁₉

106-35° (semicarbazone, m. 204-6°;

2,4-dinitrophenylhydrazone, m. 151-3°), was obtained by treating 5

g. of II with 10 cc. concentrated H₂SO₄. A small amount of

Δ 1(9)-10-methyl-2-octalone was obtained on chromatographing the

mother liquor from the hydrazone on alumina. Similarly on treatment with

I, cyclohexanone gave 2-(γ -chlorocrotyl)cyclohexanone, which

cyclized to 45% Δ 1(9)-2-octalone and

4-methylbicyclo[3.3.1]non-3-en-9-one, b₁₅ 110-15°; semicarbazone, m.

215-17°; 2,4-dinitrophenylhydrazone, m. 199-201°. Similarly

isophorone (III) and I gave 3,5,5-trimethyl-2-(γ -

chlorocrotyl)cyclohex-2-en-1-one, b₁₈ 159°, n_{15D} 1.5095;

semicarbazone, 134-6°; 2,4-dinitrophenylhydrazone, m.

133-5°, which on treatment with H₂SO₄ gave

3,5,5-trimethyl-2-(3-oxobutyl)cyclohex-2-en-1-one (IV), b_{0.1} 102°,

n_{15D} 1.4924; disemicarbazone, m. 199-201°;

mono-2,4-dinitrophenylhydrazone, m. 170-2°. IV on ozonization gave

3,3-dimethyl-5-oxohexanoic acid (V), the same product obtained by KMnO₄

oxidation of III. Condensation of III with CH₂:CHCN in the presence of Na

tert-amylate gave 3,5,5-trimethyl-2-(γ -cyanoethyl)cyclohex-2-en-1-

one (VI), b_{0.5} 124-5°, n_{21D} 1.4930; semicarbazone, m.

189-92°; 2,4-dinitrophenylhydrazone, m. 162-4°. VI was also

ozonized to V. VI on saponification gave the corresponding acid,

3,5,5-trimethyl-2-(γ -carboxyethyl)cyclohex-2-en-1-one (VII), m.

74-6°; semicarbazone, m. 211-13°. Condensation of III with

Me acrylate gave the Me ester of VII, b₁₅ 166-7°;

2,4-dinitrophenylhydrazone, m. 124-6°. Saponification of the ester gave a

substance m. 144-7, probably a lactone. IV on treatment with NaOMe in

MeOH gave 5,7,7-trimethyl-2-oxo-2,3,4,6,7,8-hexahydronaphthalene (VIII),

b_{0.1} 105°, n_{19D} 1.5542; semicarbazone, m. 206-10°;

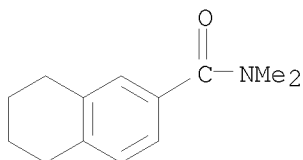
2,4-dinitrophenylhydrazone, m. 170-2°. VIII was hydrogenated over

PtO₂ to 5,7,7-trimethyl-2-decalol, m. 133-5°, which was in turn

oxidized by the method of K. Bowden, et al. (C.A. 40, 2787.8), to the

corresponding ketone; 2,4-dinitrophenylhydrazone, m. 156-9°. Dihydroisophorone and I gave 3,5,5-trimethyl-2-(γ -chlorocrotyl cyclohexanone, b15 149°, n20D 1.4851 (semicarbazone, m. 136-8°; 2,4-dinitrophenylhydrazone, m. 110-12°), which on treatment with H₂SO₄ yielded Δ 1(9)-5,7,7-trimethyl-2-octalone, b18 154°, n22D 1.5088 (semicarbazone, m. 205-8°; 2,4-dinitrophenylhydrazone, m. 184-5°). 3-Methyl-2-cyclohexen-1-one also condensed with I to form 3-methyl-2-(γ -chlorocrotyl)cyclohex-2-en-1-one, b13 146°, n21D 1.5242 (semicarbazone, m. 194-7°; 2,4-dinitrophenylhydrazone, m. 123-4°), which on treatment with H₂SO₄ yielded 3-methyl-2-(3-oxobutyl)cyclohex-2-en-1-one (IX), b14 153-4°, n18D 1.5052. IX on refluxing with Na in MeOH gave 5-methyl-2-oxo-2,3,4,6,7,8-hexahydronaphthalene, b14 156-8°, n19D 1.5738; semicarbazone, m. 195-7°; 2,4-dinitrophenylhydrazone, m. 191-4°. 4-Carbethoxy-3-methylcyclohex-2-en-1-one and I gave 4-carbethoxy-3-methyl-2-(γ -chlorocrotyl)cyclohex-2-en-1-one (X), b0.8 151-3°; semicarbazone, m. 142-4°; 2,4-dinitrophenylhydrazone, m. 122-3°. X was treated with H₂SO₄ and the neutral fraction heated with EtONa to obtain 5-methyl-2-oxo-6-carboxy-2,3,4,6,7,8-hexahydronaphthalene, m. 124-6°; 2,4-dinitrophenylhydrazone of the Me ester, m. 142-5°. (47 references.)

IT 872797-91-6P, 2-Naphthamide, 5,6,7,8-tetrahydro-N,N-dimethyl-
 RL: PREP (Preparation)
 (preparation of)
 RN 872797-91-6 CAPLUS
 CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N,N-dimethyl- (CA INDEX NAME)



L5 ANSWER 27 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1955:56623 CAPLUS

DOCUMENT NUMBER: 49:56623

ORIGINAL REFERENCE NO.: 49:10906c-e

TITLE: Syntheses and pharmacological action of Tetralin derivatives

AUTHOR(S): Fujimura, I. Hajime; Ueshima, Takaji; Fuijisawa, Toshikazu; Sugii, Michiyasu; Yaze, Toru

CORPORATE SOURCE: Univ. Kyoto

SOURCE: Yakugaku Zasshi (1954), 74, 954-6

CODEN: YKKZAJ; ISSN: 0031-6903

DOCUMENT TYPE: Journal

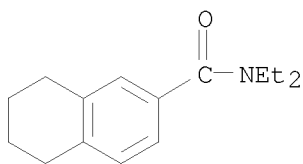
LANGUAGE: Unavailable

AB The following Tetralin derivs. are prepared: R1CH2COC1 (R1 = 5,6,7,8-tetrahydro-1-naphthyl), b9 141-2°; R2CH2COC1 (R2 = 5,6,7,8-tetrahydro-2-naphthyl), b9 144-5°. R1CH2CONR2 (R given): H, m. 135-7°; Me, b8.5 166-8°, m. 96-8°; Et, b8.5 194-6°. R1CH2CO2CH2CH2NR2.HCl (R given): Me, m. 111-13°; Et, m. 99-101°; R2CH2CONR2 (R given): H, m. 138-40°; Me, b5.5 168-70°; Et, b4.5 167-70°; R2CH2CO2CH2CH2NR2.HCl (R given): Me, m. 88-90°; Et, m. 78-80°. R2CONR2 (R given): H, m. 128-31°; Me, b6.5 163-5°; Et, b6.5 167-70°, m. 67-9°. R2CO2CH2CH2NR2.HCl (R given): Me, m. 123-6°; Et, m. 99-102°. Curarimetic, analgesic, local anesthetic, and temperature depressing actions of these products are given, although such actions are not great.

IT 856056-80-9P, 2-Naphthamide, N,N-diethyl-5,6,7,8-tetrahydro-
858199-34-5P, 2-Naphthaleneacetamide,
5,6,7,8-tetrahydro-N,N-dimethyl- 858459-67-3P,
2-Naphthaleneacetamide, N,N-diethyl-5,6,7,8-tetrahydro-
872797-91-6P, 2-Naphthamide, 5,6,7,8-tetrahydro-N,N-dimethyl-
RL: PREP (Preparation)
(preparation of)

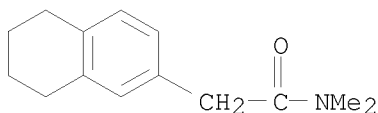
RN 856056-80-9 CAPLUS

CN 2-Naphthalenecarboxamide, N,N-diethyl-5,6,7,8-tetrahydro- (CA INDEX NAME)



RN 858199-34-5 CAPLUS

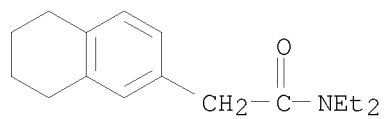
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N,N-dimethyl- (CA INDEX NAME)



RN 858459-67-3 CAPLUS

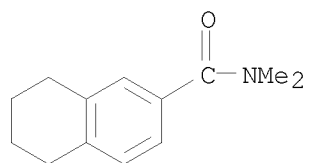
CN 2-Naphthaleneacetamide, N,N-diethyl-5,6,7,8-tetrahydro- (CA INDEX NAME)

10/513699



RN 872797-91-6 CAPLUS

CN 2-Naphthalenecarboxamide, 5,6,7,8-tetrahydro-N,N-dimethyl- (CA INDEX NAME)



<12/04/2007>

Erich Leese

L5 ANSWER 28 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1954:61453 CAPLUS

DOCUMENT NUMBER: 48:61453

ORIGINAL REFERENCE NO.: 48:10920c-e

TITLE: Further investigation on the mitosis-poison action of dihydrostilbylamine derivatives

AUTHOR(S): Lettre, Hans; Delitzsch, Ingrid

CORPORATE SOURCE: Univ. Gottingen, Germany

SOURCE: Hoppe-Seyler's Zeitschrift fuer Physiologische Chemie (1952), 289, 220-5

CODEN: HSZPAZ; ISSN: 0018-4888

DOCUMENT TYPE: Journal

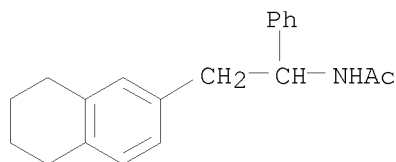
LANGUAGE: Unavailable

AB cf. C.A. 37, 5784.8(1943). Phenylnitromethane and 3,4-ethylenedioxybenzaldehyde are condensed in the presence of $\text{CH}_3\text{NH}_2\text{-HCl}$ to give 3',4'-ethylenedioxy- α -nitrostilbene, m.p. 142-3°; this compound is reduced with Zn dust and AcOH and subsequently Na-Hg to give the corresponding dihydrostilbylamine-HCl, m.p. 225-8°. Its activity against in vitro fibroblasts (0.75 γ /ml.) is 4 times that of the corresponding dioxymethylene compound (3 γ /ml.). N-Methylformanilide and POCl_3 are treated with catechol tetramethylene ether; the aldehyde obtained by this method is condensed with phenylnitromethane and reduced to give 3',4'-tetramethylenedioxydihydrostilbylamine. This compound shows no mitosis-poison activity. The 4'-methoxy derivative shows activity at 4-5 γ /ml., 4'-methylstilbylamine at 10 γ /ml., 3',4'-trimethylenedioxydihydrostilbylamine at 10 γ /ml. The two latter compds. are prepared from the formyl derivs. of hydrindene and Tetralin, resp. No mitosis-poison effect was shown by the corresponding aromatic compds. made from 1- and 2-naphthaldehydes. This is an opposite trend as observed for the carcinogenic properties of benzopyrene, which is very active compared with its tetrahydro derivative

IT 855928-78-8P, Acetamide,
N-[α -(5,6,7,8-tetrahydro-2-naphthyl)methyl]benzyl]-
858459-72-0P, Naphthalene,
1,2,3,4-tetrahydro-6-(β -nitrostyryl)-
RL: PREP (Preparation)
(preparation of)

RN 855928-78-8 CAPLUS

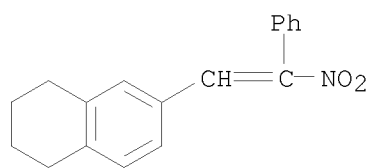
CN Acetamide, N-[1-phenyl-2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]- (CA INDEX NAME)



RN 858459-72-0 CAPLUS

CN Naphthalene, 1,2,3,4-tetrahydro-6-(2-nitro-2-phenylethenyl)- (CA INDEX NAME)

10/513699



<12/04/2007>

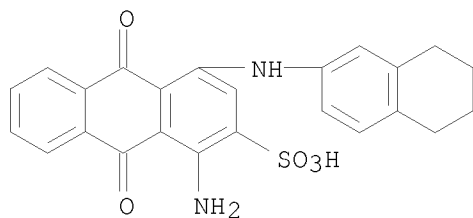
Erich Leese

10/513699

L5 ANSWER 29 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1954:48257 CAPLUS
DOCUMENT NUMBER: 48:48257
ORIGINAL REFERENCE NO.: 48:8551i,8552a-b
TITLE: Acid fulling dyes of the anthraquinone series
PATENT ASSIGNEE(S): Sandoz Ltd.
DOCUMENT TYPE: Patent
LANGUAGE: Unavailable
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
	-----	----	-----	-----	-----
	GB 699781		19531118	GB 1951-19700	19510821 <--
AB	Compds. having the formula 1,2,4-H ₂ N(MO ₃ S)(RHN)C ₁₄ H ₅ O ₂ (I), where M is H, NH ₄ , Li, Na, or K and R is substituted or unsubstituted tetrahydronaphthyl or C ₆ H ₄ NHCOPh, are brominated to give acid fulling dyes. Thus, Na 1-amino-4-(1,2,3,4-tetrahydro-5-naphthylamino)-2-anthraquinonesulfonate 47 (obtained by condensing the Na 1-amino-4-bromo-2-anthraquinonesulfonate in H ₂ O with Cu as a catalyst with 5-amino-1,2,3,4-tetrahydronaphthalene) is dissolved in 90% H ₂ SO ₄ 1500, treated with Br 26, and the mixture stirred for 6 hrs. at room temperature and then 3 hrs. at 60° to give a brominated dye (II). I dyes wool and other animal fibers as well as nylon with a red-tinged blue tint. Other dyes are similarly obtained by brominating I, where R = 1,2,3,4-tetrahydro-6-naphthyl, p-PhCONHC ₆ H ₄ , p-(p-NH ₂ C ₆ H ₄ CONH)C ₆ H ₄ , p-[2,5-Cl(AcNH)C ₆ H ₃ CONH]C ₆ H ₄ , p-[3,4-Me(H ₂ N)C ₆ H ₃ CONH]C ₆ H ₄ , and p-(p-FC ₆ H ₄ CONH)C ₆ H ₄ .				
IT	859335-68-5, 2-Anthraquinonesulfonic acid, 1-amino-4-[5,6,7,8-tetrahydro-2-naphthylamino]- (and bromine derivs.)				
RN	859335-68-5 CAPLUS				
CN	2-Anthracenesulfonic acid, 1-amino-9,10-dihydro-9,10-dioxo-4-[(5,6,7,8-tetrahydro-2-naphthalenyl)amino]- (CA INDEX NAME)				



L5 ANSWER 30 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1952:54594 CAPLUS

DOCUMENT NUMBER: 46:54594

ORIGINAL REFERENCE NO.: 46:9092i,9093a-i

TITLE: Carcinogenic nitrogen compounds. IX. The use of aminotetralins for the synthesis of dibenzacridines and related compounds

AUTHOR(S): Buu-Hoi, Ng. Ph.; Jacquignon, Pierre

CORPORATE SOURCE: Univ. Paris

SOURCE: Journal of the Chemical Society (1951)
2964-8

CODEN: JCSEA9; ISSN: 0368-1769

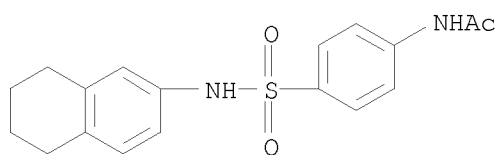
DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

AB 5,6,7,8-Tetrahydro-2-naphthylamine (I) (preparation given) yields a p-toluenesulfonate, m. 137°, and a N-(N-acetylsulfanilyl) derivative, m. 216°. 6-Acetyl-1,2,3,4-tetrahydronaphthalene (100 g.) on Huang-Minlon reduction gives 70 g. 6-ethyl-1,2,3,4-tetrahydronaphthalene (II), b. 239°; with AcCl and AlCl₃ in CS₂ II yields 6-acetyl-7-ethyl-1,2,3,4-tetrahydronaphthalene, pale yellow, b₁₇, 188°; semicarbazone, m. 134-5°; thiosemicarbazone, m. 123°; oxime (III), m. 128°. III (33 g.), and 30 g. PCl₅ in 80 cc. ether give 32.5 g. of the N-Ac derivative, m. 166°, of 3-ethyl-5,6,7,8-tetrahydro-2-naphthylamine (IV), pale yellow, b₁₆ 175°; HCl salt, m. 147°; p-toluenesulfonate, m. 130°; N-(N-acetylsulfanilyl) derivative, m. 201°. I (5 g.), 5 g. 1-C₁₀H₇OH, and 0.75 g. (HCHO)₃ give 2 g. 1'',2'',3'',4''-tetrahydro-1,2:7,8-dibenzacridine (V), pale yellow, m. 118° [picrate, orange-red, m. 257° (decomposition)]; 0.5 g. V and 0.3 g. Se, heated 3 h. at 350°, give 1,2:7,8-dibenzacridine, pale yellow, m. 129° [picrate, brick-red, m. 266° (decomposition)]; 2-C₁₀H₇OH gives 4 g. 1',2',3',4'-tetrahydro-2,3:6,7-dibenzacridine (VI), pale yellow, m. 145°. 6,2-(tert-Bu)C₁₀H₆OH (2.2 g.) gives about 3 g. of the 3'-tert-Bu derivative of VI, pale yellow, m. 128° (picrate, orange, decompose above 231°). I (7.5 g.), 12.5 g. 1-C₁₀H₇OH, and 0.1 g. iodine, heated 5 h. at 240-5°, give 12 g. N-(5,6,7,8-tetrahydro-2-naphthyl)-1-naphthylamine (VII), yellow, b₁₆ 285-90°; 2-C₁₀H₇OH yields the 2-isomer (VIII), b₁₆ 304-5°, m. 96°. VII (2 g.), 2 g. Ac₂O, and 2 g. ZnCl₂, heated 6 h. at 180-90°, and the product treated with hot aqueous NaOH and extracted with PhMe, give 1.5 g. 1'',2'',3'',4''-tetrahydro-5-methyl-1,2:7,8-dibenzacridine, pale yellow, m. 130° (picrate, orange-yellow, m. 251°); (EtCO)₂O gives the 5-Et homolog, yellow, m. 114° (picrate, orange-yellow, m. 223°). VIII (5 g.), Ac₂O, and ZnCl₂ give 3 g. 1',2',3',4'-tetrahydro-5-methyl-2,3:6,7-dibenzacridine, pale yellow, m. 166° (picrate, brownish red, m. 272-3°); 5-Et homolog, pale yellow, m. 171° (picrate, bright red, m. 253°). VII (4 g.) and 2 g. AsCl₃ in 20 cc. o-C₆H₄Cl₂, refluxed 2 h., give 3.7 g. 10-chloro-1',2',3',4',5,10-hexahydro-2,3:6,7-dibenzophenarsazine, orange-yellow, m. 264°; with MeMgI this yields 1',2',3',4',5,10-hexahydro-10-methyl-2,3:6,7-dibenzophenarsazine, m. 226°; 10-Et homolog, m. 157°. VIII (4 g.) yields 3.7 g. 1'',2'',3'',4'',5,10-hexahydro-2,3:6,7-dibenzophenarsazine, orange yellow, m. 260°; the 10-M derivative, m. 200°, and the 10-Et homolog, m. 146°. IV (4 g.), 4 g. 1-C₁₀H₇OH, and 0.75 g. (HCHO)₃ give 9-ethyl-1'',2'',3'',4''-tetrahydro-1,2:6,7-dibenzacridine, pale greenish yellow, m. 132° (picrate, orange-red, m. 173°); heated with

Se (3 h. at 350°), this yields 9-ethyl-1,2:6,7-dibenzacridine, pale yellow, m. 110° (picrate, light orange, m. 256°). IV and 2-C10H7OH with (HCHO)3 give 1-ethyl-1',2',3',4'-tetrahydro-3,4:6,7-dibenzacridine (IX), yellow, m. 140° (picrate, orange, m. 298°); 0.5 g. IX and 0.3 g. Se give 0.3 g. 1-ethyl-3,4:6,7-dibenzacridine, pale yellow, m. 158° (picrate, yellow, m. 276°). 6,2-(tert-Bu)C10H6OH, IV, and (HCHO)3 give 3''-tert-butyl-1',2',3',4'-tetrahydro(3,4:6,7)dibenzacridine, yellow, m. 154° (picrate, orange yellow, m. 257-8°). I (5 g.), 5 g. Ac2CH2, and 3 drops AcOH, heated 16 h. at 170-80° give 5 g. 2,5-dimethyl-1-(5,6,7,8-tetrahydro-2-naphthyl)pyrrole, pale yellow, b16 200-2°, nD21.5 1.5790; IV yields 5 g. of the 1-(3-ethyl-5,6,7,8-tetrahydro-2-naphthyl) homolog, pale yellow, b17, 215-16°. IV (2 g.) and 2 g. Ac2CH2, refluxed 2 h., and the cold solution treated with 12 cc. H2SO4 and heated 1 h. on the water bath, give 1.8 g. 8-ethyl-1',2',3',4'-tetrahydro-2,4-dimethyl-5,6-benzoquinoline, pale yellow, b17, 214-16°, nD21.5 1.6125 (picrate, yellow, m. 223-4°). 5,6,7,8-Tetrahydro-2-naphthylhydrazine-HCl (1 g.), 1 g. 1-indanone, and 1 g. AcONa in 20 cc. EtOH, refluxed 1 h. and the crude hydrazone heated a few sec. with HCl in AcOH, give 1 g. 1'',2'',3'',4''-tetrahydro-5,6-benzindeno(3',2':2,3)indole, m. 297°; 1 g. 3,4-dihydro-1(2H)-naphthalenone yields 1.2 g. 1'',2'',3,3'',4,4''-hexahydro-1,2:6,7-dibenzocarbazole, m. 190° (picrate, brown-violet, m. 173°); 1 g. 1-oxo-1,2,3,4,5,6,7,8-octahydroanthracene gives 1.5 g. 1'',2'',3,3'',4,4'',5',6',7',8'-decahydro-6,7-benzonaphtho(2',3':1,2)carbazole, m. 208° (picrate, deep violet, m. 203°).

IT 857552-84-2P, Acetanilide,
4'-[(5,6,7,8-tetrahydro-2-naphthyl)sulfamoyl]-
RL: PREP (Preparation)
(preparation of)
RN 857552-84-2 CAPLUS
CN Acetamide, N-[4-[(5,6,7,8-tetrahydro-2-naphthalenyl)amino]sulfonyl]phenyl]- (CA INDEX NAME)



OS.CITING REF COUNT: 1 THERE ARE 1 CAPLUS RECORDS THAT CITE THIS RECORD
(1 CITINGS)

L5 ANSWER 31 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1950:10099 CAPLUS
 DOCUMENT NUMBER: 44:10099
 ORIGINAL REFERENCE NO.: 44:1981i,1982a-i,1983a-e
 TITLE: Orienting phenomena in the substitution on aromatic bicyclic nuclei. II. Combe's quinoline synthesis
 AUTHOR(S): Huisgen, Rolf
 SOURCE: Justus Liebig's Annalen der Chemie (1949), 564, 16-32
 CODEN: JLACBF; ISSN: 0075-4617
 DOCUMENT TYPE: Journal
 LANGUAGE: Unavailable
 OTHER SOURCE(S): CASREACT 44:10099

GI For diagram(s), see printed CA Issue.

AB cf. C.A. 42, 6783i. When Combe's synthesis [Compt. rend. 106, 142, 1536 (1888)] was applied to 2-C10H7NH2 (I) by Johnson and Mathews (C.A. 38, 1505.6), only linear cyclization was effected, which is an exception to the usually observed angular cyclization reaction. The structure of the side chain in position 2 on the C10H8 nucleus is evidently important, and the usually accepted -N:CRCH2COR appears less satisfactory than does the "enamine" structure, -NHCR:CHCOR, which is adopted by H. Conditions for forming nonanellated products and the theory underlying linear ring closure are discussed at length. 2-(Ac.CH:CMcNH)C10H7 (II), m. 99°, was prepared by heating I with Ac2CH2 (III). On oxidation, II gave the Ac derivative of I, m. 132-3°, and, when heated at 290° in paraffin oil, followed by extraction with MeOH, evaporation, extraction with C6H6, and acidification, gave I.HCl. At 0° 1 g. II and 6 cc. concentrated H2SO4 gave 2,4-dimethylbenzo[g]quinoline (IV), but with H3PO4 at 70°, followed by extraction with NH3 in CHCl3, II gave largely I and only small amts. of IV. By the cyclization of appropriate enamines the following derivs. of IV were formed: 8-Br, pale yellow, m. 168-9° (from the enamine prepared from III and 1,2-H2N(Br)C10H6), and 8-Me, m. 127° (prepared from the enamine, m. 95°, formed from III and 2-H2NC10H6Me). The 3-Br derivative (V) of I and III gave 2-(3-bromo-2-naphthylamino)-2-penten-4-one, needles, m. 95° (from Et2O), which with concentrated H2SO4 at 0° gave 83% V. I (10 g.) and 17 g. Bz2CH2 at 130° gave the enamine, 2-(BzCH:CPhNH)-C10H7 (VI), yellow needles, m. 146-7°, nonfluorescent, unaffected by hot aqueous acids or alkalies, and unchanged when heated at 300° in paraffin, but forming 2-BzNHC10H7, m. 159-60° and BzOK when oxidized with KMnO4 in Me2CO. With 90 cc. concentrated H2SO4 at 0°, 15 g. VI gave an orange-red sulfate which was washed with ice-H2O, dried, extracted with H2O and CHCl3, and the dried CHCl3 extract was passed through Al2O3, evaporated, and crystallized from EtOH at 40°, giving 10.81 g. 2,4-diphenylbenzo[g]quinoline (VII), pale yellow, m. 146° (from CHCl3-EtOH), showing in alc. an unusually brilliant pale blue fluorescence veering to faint yellow on addition of acid, and, from the mother liquor of VII, 1.2 g. of the angular [f]-isomer (VIII), m. 146°, separated after tedious, successive fractionations from Me2CO, AcOEt, and MeOH. A simpler means of separating VII and VIII was to irradiate the mixture in Me2CO, thus forming the insol. dimer (IX) of VII, the mother liquors from which gave nearly pure VIII. When mixed, VII and VIII showed m.-p. depressions of 20-30°. With glacial AcOH and CrO3, 1.5 g. VII formed an insol. orange-red bichromate, which, when heated, gave 1.39 g. yellow 2,4-diphenyl-1-azanthraquinone, C25H15NO2, m. 250-1° (from glacial

AcOH or C₆H₆). VII (0.3 g.) gave 0.375 g. of a sulfonic acid, C₂₅H₁₆NSO₃H, orange-red needles (from EtOH), m. above 300°. Heated with quinone in C₆H₆ VII gave the adduct (X), colorless, nonfluorescent, dissociating and foaming at 204-6°, and giving a clear melt at 240°. A 24-hr. solar photodimerization of VII in C₆H₆ gave IX, C₅₀H₃₁N₂, colorless, m. 273 ° (from C₆H₆), showing a faint blue fluorescence and quantitatively depolymerized to VII on melting. VIII was effectively prepared by gradually adding 0.5 g. VI to ZnCl₂ at 200° heating 1 hr., decomposing the melt with H₂O, extracting with CHCl₃, passing the

extract

through Al₂O₃, evaporating, and crystg, from MeOH; yield 0.28 g. VIII. VIII in alc. showed a bluish-violet fluorescence; VIII forms difficultly soluble HCl and H₂SO₄ salts. In alc., VIII proved stable on irradiation, but a photochem. reaction occurred rapidly on addition of a few drops of H₂SO₄, giving (from 0.15 g. VIII) 0.136 g. of a compound (XI), felted needles, m. 220° (from CHCl₃-EtOH). Bz₂CH₂ and the 1-Me derivative of I gave the corresponding enamine, m. 163-4°, which when cyclized at 5° with H₂SO₄ gave 96% of the 8-Me derivative (XII) of VII, m. 138-9° (from EtOH-CHCl₃), whose photodimer, colorless crystals with bluish-green surface sheen, m. 246-7° XII when added to a ZnCl₂ melt gave a compound, (1-MeC₁₀H₆)₂NH (?), m. 221-4° (insol. even in concentrated acids). The 1-Br derivative of I and Bz₂CH₂ gave an enamine, yellow leaflets, m. 174°, difficultly cyclized after standing 24 hrs. in concentrated H₂SO₄ to give a poor yield of an impure (linear ?) pale yellow bromobenzoquinoline, m. 199-204° (containing 77.41% C instead of the calculated 73.16%); when this enamine was heated with ZnCl₂, small amts. of VIII were formed. The 3-Br derivative of I and Bz₂CH₂ reacted slowly at 130° to give the enamine, C₂₅H₁₈NOBr, m. 131°, which when fused with ZnCl₂ gave 1,3-diphenyl-5-bromobenzo[f]quinoline, m. 152-3° showing weak violet fluorescence. Bz₂CH₂ and freshly distilled 2-amino-1,2,3,4-tetrahydronaphthalene gave an enamine, m. 135° which with H₂SO₄ at 5° gave 92% (linear) 2,4-diphenyl-6,7,8,9-tetrahydrobenzo[g]quinoline (XIII), m. 129°, showing a blue fluorescence, also formed from the enamine by ZnCl₂ fusion. Dehydrogenation of XIII with Pt-C gave VII; no VIII was formed. I and BzCH₂Ac at 130° gave 96% of the corresponding enamine, m. 152-3°, which yielded 2-methyl-4-phenylbenzo[g]quinoline (XIV), b₁₂ 260°, m. 110° (from petr. ether) [cf. Beyer, Ber. 20, 1767 (1887)]. Similarly, the 1-Me derivative of I and BzCH₂Ac gave 90% of an enamine, pale yellow prisms, m. 158°, which gave 70% of the 8-Me deriv, of XIV, m. 105-6°. The enamine, m. 133-4°, prepared from 3,2-BrC₁₀H₆-NH₂ and BzCH₂Ac could not be cyclized by the usual procedure with H₂SO₄. Ultraviolet absorption spectra of VII and VIII are given.

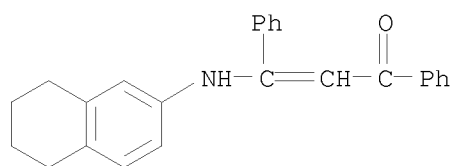
IT 854835-41-9P, Chalcone, β-(5,6,7,8-tetrahydro-2-naphthylamino)-

RL: PREP (Preparation)
(preparation of)

RN 854835-41-9 CAPLUS

CN 2-Propen-1-one, 1,3-diphenyl-3-[(5,6,7,8-tetrahydro-2-naphthalenyl)amino]-
(CA INDEX NAME)

10/513699



OS.CITING REF COUNT: 2

THERE ARE 2 CAPLUS RECORDS THAT CITE THIS RECORD
(2 CITINGS)

L5 ANSWER 32 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1949:46455 CAPLUS

DOCUMENT NUMBER: 43:46455

ORIGINAL REFERENCE NO.: 43:8390a-i

TITLE: Synthetic studies in the isoquinoline series

AUTHOR(S): Schultz, Everett M.; Arnold, R. T.

SOURCE: Journal of the American Chemical Society (1949), 71, 1911-14

CODEN: JACSAT; ISSN: 0002-7863

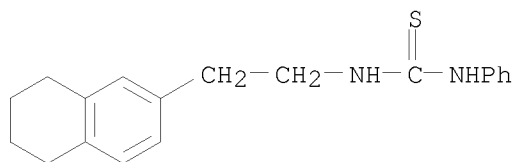
DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

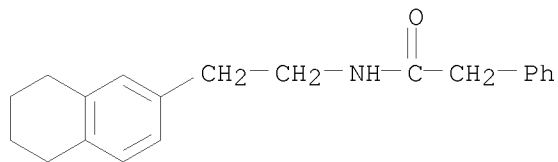
AB 5-(Chloromethyl)hydrindene (30 g.) in 75 ml. EtOH, added (0.5 hr.) to 11 g. NaCN in a min. quantity of hot H₂O and refluxed 4 hrs., gives 73% 5-hydrindeneacetonitrile (I), b_{2.5} 113-15°; 58 g. I and 4 g. Raney Ni in 100 ml. EtOH containing 19 g. NH₃, hydrogenated (4 hrs.) at 110° and 1100 lb., give 98% 2-(5-hydrindenyl)ethylamine (II), b₄ 104-7°, analyzed as the phenylthiocarbamyl derivative, m. 96.5-7.5°. I (75 g.) in 24 g. absolute EtOH and 70 ml. ether, treated with 21 g. HCl, kept 16 hrs. at 5°, and the imido ester hydrolyzed (24 hrs.) with H₂O, gives 40 g. of the Et ester, b₃ 122°, n_D²⁵ 1.5201, of 5-hydrindeneacetic acid (III), m. 113-14° (anilide, m. 122.5-3.5°). The acid chloride of III (19.4 g.), added to 16.1 g. II in 7.6 g. C₅H₅N and boiled 5 min., gives 47% N-[2-(5-hydrindenyl)ethyl]-5-hydrindeneacetamide (IV), m. 99-100°. IV (6 g.) in 65 ml. hot xylene, added to 12 g. P₂O₅, the mixture boiled 15 min., 6 g. P₂O₅ added, the mixture boiled an addnl. 0.5 hr., the xylene decanted, the residue heated with 100 ml. H₂O, the residual xylene removed by steam distillation, and the aqueous solution extracted with ether, made strongly basic with 10% KOH, and extracted with ether, gives 4 g. 1-(5-hydrindenylmethyl)-6,7-cyclopenteno-3,4-dihydroisoquinoline (V), yellow, b₃ 235-40° (bath temperature), analyzed as the picrate, m. 196-7°; the crude V and 0.6 g. 10% Pd-C, heated in a CO₂ stream 45 min. at 180-200°, give 1.3 g. 1-(5-hydrindenylmethyl)-6,7-cyclopentenoisoquinoline, m. 91-2°. II (8.86 g.) and 11 g. homopiperonylic acid, heated 4 hrs. at 160-70°, give 13.9 g. N-[2-(5-hydrindenyl)ethyl]homopiperonylamide (VI), m. 119-20°; 5.5 g. VI, cyclized as above and dehydrogenated, gives 0.8 g. 1-(3,4-methylenedioxybenzyl)-6,7-cyclopentenoisoquinoline, m. 98-9°. The crude acid chloride from 17.6 g. III in 100 ml. C₆H₆, treated slowly with 16.5 g. homopiperonylamine and 7.6 g. C₅H₅N in 50 ml. C₆H₆ and boiled 10 min., give 61% N-[2-(3,4-methylenedioxyphenyl)-ethyl]-5-hydrindeneacetamide, m. 122.5-3.5°; cyclization of 2 g. with P₂O₅ gives 0.7-0.8 g. 1-(5-hydrindenylmethyl)-6,7-methylenedioxy-3,4-dihydroisoquinoline (VIa), m. 130.5-1° (picrate, m. 175-6°); 5.9 g. (VIa) and 1.3 g. 10% Pd-C, heated 3.5 hrs. at 155-200°, give 1.65 g. of the HCl salt, m. 257-8°, of 1-(5-hydrindenylmethyl)-6,7-methylenedioxyisoquinoline, m. 168-9° (picrate, m. 184-5°). II (5.33 g.), 2.57 g. AcCl, and 2.6 g. C₅H₅N in C₆H₆ give 3 g. of the Ac derivative, b₂ 170-85° (bath temperature), m. 77.5-8°; cyclization gives 1-methyl-6,7-cyclopenteno-3,4-dihydroisoquinoline, analyzed as the picrate, m. 205°; oxidation of the base with HNO₃ gives 1,2,4,5-C₆H₂(CO₂H)₄, indicating that cyclization in the hydrindene series occurs across the 5,6-positions of the hydrindene nucleus. 6-Acetyl-1,2,3,4-tetrahydronaphthalene (50.9 g.), 10 g. S, and 26 g. morpholine, heated 8.5 hrs. at 120-5°, give 90% 1,2,3,4-tetrahydro-6-naphthalenethioacetomorpholide (VII), m. 114.5-15.5°; 58.4 g. VII in 1 l. 10% KOH, boiled 10 hrs., gives 79%

1,2,3,4-tetrahydro-6-naphthaleneacetic acid (VIII), m. 95-6°; the Me ester of VIII (b2 141-5°) and concentrated NH₄OH, 72 hrs. at 25-30°, give 83% of the amide, m. 168-9°; with SOCl₂ in C₆H₆ this yields 67% of the nitrile, b3 144-7°, catalytic reduction of which in MeOH containing liquid NH₃ gives 90% 2-(1,2,3,4-tetrahydro-6-naphthyl)ethylamine (IX), analyzed as 1-[2-(1,2,3,4-tetrahydro-6-naphthyl)ethyl]-3-phenyl-2-thiourea, m. 130-1°; 5 g. IX and 4.25 g. PhCH₂CO₂H, heated 3 hrs. at 160-80°, give 74% N-[2-(1,2,3,4-tetrahydro-6-naphthyl)ethyl]-α-phenylacetamide, m. 99-100°; cyclization of 5.15 g. with 10 g. P₂O₅ in 75 ml. PhMe gives 3.75 g. 1-benzyl-6,7-cyclohexeno-3,4-dihydroisoquinoline, b. 180-200°/10-4 mm., analyzed as the picrate, m. 193-4° (decomposition); HNO₃ oxidation yields 1,2,4,5-C₆H₂(CO₂H)₄; heating with 10% Pd-C at 300-310° (4.5 hrs.) gives the isoquinoline (m. 115-16°), whose picrate m. 211-12°.

IT 859736-98-4P, Urea, 1-phenyl-3-[2-(5,6,7,8-tetrahydro-2-naphthyl)ethyl]-2-thio- 861059-29-2P, Acetamide, 2-phenyl-N-[2-(5,6,7,8-tetrahydro-2-naphthyl)ethyl]-
 RL: PREP (Preparation)
 (preparation of)
 RN 859736-98-4 CAPLUS
 CN Thiourea, N-phenyl-N'-[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]- (CA INDEX NAME)



RN 861059-29-2 CAPLUS
 CN Benzeneacetamide, N-[2-(5,6,7,8-tetrahydro-2-naphthalenyl)ethyl]- (CA INDEX NAME)



OS.CITING REF COUNT: 1 THERE ARE 1 CAPLUS RECORDS THAT CITE THIS RECORD
 (1 CITINGS)

L5 ANSWER 33 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1949:17411 CAPLUS

DOCUMENT NUMBER: 43:17411

ORIGINAL REFERENCE NO.: 43:3362g-i,3363a-i,3364a-d

TITLE: Biosynthesis of penicillins. VI. N-2-Hydroxyethyl amides of some polycyclic and heterocyclic acetic acids as precursors

AUTHOR(S): Jones, Reuben G.; Soper, Quentin F.; Behrens, Otto K.; Corse, Joseph W.

SOURCE: Journal of the American Chemical Society (1948), 70, 2843-8

CODEN: JACSAT; ISSN: 0002-7863

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

OTHER SOURCE(S): CASREACT 43:17411

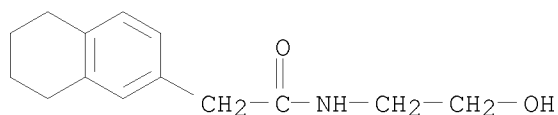
AB 2,6-MeCl₁₀H₆NH₂ (78 g.) in 80 mL. concentrated HCl and 200 mL. H₂O at 0°, treated at 5° with 35 g. NaNO₂ in 50 mL. H₂O and, after 0.5 h., with 130 g. ice-cold 42% HBF₄, gives 90% of the 2-diazonium fluoroborate, decomposition of which yields 69% 2-methyl-6-fluoronaphthalene (I), m. 77°. I (40 g.) at 210°, treated (15 min.) with 40 g. Br (with illumination with a 100-w. lamp), gives 82% 2-(bromomethyl)-6-fluoronaphthalene (II), b₂ 125-30°, m. 53°. II (48 g.), added to a refluxing solution of 30 g. KCN in 60 mL. H₂O and 200 mL. EtOH, the EtOH removed after refluxing 4 h., 500 mL. H₂O added, the solution extracted with ether, and the residue from the ether boiled 5 h. with 40 g. KOH in 40 mL. H₂O and 200 mL. EtOH, gives 74% 6-fluoro-2-naphthaleneacetic acid, m. 138-9° (Me ester, b₂ 163-6°, m. 48-9°). 2,6-MeCl₁₀H₆NH₂ (63 g.) in 100 mL. H₂O and 700 g. 48% HBr, treated (3-4 h.) at 5° with 45 g. NaNO₂ in 75 mL. H₂O and the diazonium solution poured (10 min.) into 170 g. CuBr in 800 mL. 48% HBr at 70-80°, gives 40% 6-bromo-2-methylnaphthalene (III), m. 142° III yields 80% 6-bromo-2-(bromomethyl)naphthalene, m. 124-5° this gives 69% 6-bromo-2-naphthaleneacetic acid, m. 175-6° (Me ester, b₂ 187-93°, m. 67-9°). 3,2-ClCl₁₀H₆CHO (32.5 g.), 35 g. hippuric acid, 14.5 g. anhydrous AcONa, and 50 mL. Ac₂O, heated on the steam bath 1 h., give 75% 2-phenyl-4-(3-chloro-2-naphthylmethylene)-5(4H)-oxazolone (IV), bright yellow, m. 192° 40 g. IV in 200 mL. 10% NaOH, refluxed 9 h., the mixture diluted to 1500 mL. with H₂O, washed with ether, the aqueous solution treated with 20 mL. 12.5 N NaOH and 15 mL. 30% H₂O₂, allowed to stand overnight, the filtrate acidified with HCl, extracted with ether-C₆H₆, and the residue esterified, gives 37% Me 3-chloro-2-naphthaleneacetate, b₂ 163-5°, m. 49-50° the free acid m. 193-4°. 6,2-MeOC₁₀H₆Ac (100 g.), 25.5 g. S, and 87 g. morpholine, heated 18 h. at 140°, part of the morpholine removed in vacuo, 250 mL. AcOH and 350 mL. concentrated HCl added, and the mixture refluxed 24 h., give 67% 6-methoxy-2-naphthaleneacetic acid, m. 203-5° (Me ester, b₁ 192-3°, m. 86°, 73%). 5,6,7,8-Tetrahydro-2-acetonaphthone (50 g.), 13 g. S, and 40 mL. morpholine, refluxed overnight, 400 mL. concentrated HCl and 300 mL. H₂O added, and the mixture again refluxed overnight, followed by esterification with EtOH and H₂SO₄, give Et 5,6,7,8-tetrahydro-2-naphthaleneacetate, b_{0.5} 140-3°. 2-Acetylphenanthrene (13.2 g.), 3.2 g. S, and 10.5 g. morpholine, heated 15 h. at 160°, the mixture treated with 150 mL. AcOH and 36% HCl, and refluxed 24 h., give 81% 2-phenanthreneacetic acid, m. 187-8° the 3-isomer m. 174-5°, 84% (Me ester, b_{1.5}

203-5°, 89%). 8-(Bromomethyl)quinoline (120 g.) in 250 mL. warm EtOH, added (0.5 h.) to 50 g. KCN in 100 mL. warm H₂O and the mixture refluxed 1.5 h., gives 78% 8-(cyanomethyl)quinoline, m. 86-7°; hydrolysis with aqueous alc. KOH and esterification give 91% Et 8-quinolineacetate, b₃ 158-60°. Et 3-quinolinecarboxylate (70 g.), 62 g. AcOEt, and EtONa (12 g. Na and 0.52 mol absolute EtOH) in 100 cc. dry C₆H₆, refluxed 20 h., the cooled solution poured onto ice, diluted to 5 l. with H₂O, treated with 50 mL. 12 N NaOH, washed with two 300 mL. portions of ether, and the aqueous solution neutralized with dilute H₂SO₄ and extracted with two 500-mL. portions of ether, give 75% Et 3-quinolylformylacetate, m. 84° 27 g. of the keto ester in 125 g. 25% H₂SO₄, heated 30 min. at 100°, gives 95% 3-acetylquinoline (V). V (7 g.), 5 g. S, 50 mL. (NH₄)₂S, and 25 mL. H₂O, heated 20 h. at 145-50°, the residue extracted with two 300-mL. portions boiling 5% HCl, the solution refluxed 3 h., and the crude acid esterified, give 19% Et 3-quinolineacetate, b_{2.5} 140-2°. pH₂NC₆H₄CH₂CO₂H (46 g.), 10.5 g. FeSO₄, 115 g. C₃H₅(OH)₃, 23 g. PhNO₂, and 53 mL. concentrated H₂SO₄, boiled 5 h., give 37 g. crude acid which, esterified with EtOH and HCl, gives 39% Et 6-quinolineacetate, b₃ 160° the free acid (VI) m. 218-20°. Et 6-quinolinecarboxylate and AcOEt, condensed with EtONa, give 87% Et 6-quinolineacetate, hydrolysis of which with 25% H₂SO₄ at 100° gives 90% 6-acetylquinoline, m. 76° the Willgerodt reaction gives 87.5% VI. 3,4 O₂N(H₂N)C₆H₃CO₂H (108 g.) in 350 mL. concentrated HCl, treated with 125 g. Sn in portions (temperature below 90°), gives 87% (3,4-diaminophenyl)acetic acid-2HCl (VII), m. 222-4° (decomposition); Et ester-2HCl (VIII), m. 185-7° (decomposition); 3 g. VII and 20 mL. 98-100% HCO₂H, heated several hrs., give 100% 5-benzimidazoleacetic acid-HCl, m. 240-2° the Et ester m. 65-6°, 75%. VIII (14 g.) in 200 mL. ice H₂O, treated with excess COCl₂, gives 95% Et 2-hydroxy-5-benzimidazoleacetate, m. 208-9°. NCCH₂CO₂Et (113 g.) and 15 g. (HOCH₂CH₂)₃N in 100 mL. absolute EtOH, treated with a slow stream of H₂S, the mixture poured after 5 days into ice-H₂O, and 38 g. of the resulting oil and 23.1 g. ClCH₂Ac in 300 cc. anhydrous ether kept 4 days, give 20.6 g. Et 4-methyl-2-thiazoleacetate, b₁₇ 136-9°. Thiaxanthidrol (42 g.), 30 g. CH₂(CO₂H)₂, and 80 mL. C₅H₅N, heated 2 h. at 60-70° and 2 h. at 90-5° and the liquid poured into 600 mL. 2 N HCl, give 90% 9-thiaxantheneacetic acid, m. 167-8° (Me ester, b₂ 182-4°). The Ag salt of 2-benzylimidazole (53 g.) and 50 g. BrCH₂CO₂Et in 200 mL. xylene, refluxed 48 h., give 25.4% of the Et ester, m. 70-70.5°, of 2-benzyl-1-imidazoleacetic acid, m. 173-4°. Me 1-acenaphtheneacetate, b₄ 176-8°. N-2-Thienylacetyl-DL-valine m. 110-12°. Amides were prepared by heating the Me or Et ester of the various acids with a slight excess of HOCH₂CH₂NH₂ at 100-150° for several hrs.; R in RCH₂CONHCH₂CH₂OH is given, together with S (see part V). 2-Cl₁₀H₇ m. 125-7°, S 1.3; 1-bromo-2-naphthalene m. 155-6°, S 0.5; 6-fluoro-2-naphthalene m. 145-6°, S 1.2; 3-chloro-2-naphthalene m. 150-1°, S 0.3; 6-bromo-2-naphthalene m. 167-8°, S 0.9; 5,6,7,8-tetrahydro-2-naphthalene m. 88-90°, S 0.9; 1-nitro-2-naphthalene m. 154-5°, S 0.9; 6-methoxy-2-naphthalene m. 160°, S 1.1; 1-acenaphthene m. 160°, S 1.1; 9-fluorene m. 127-8°, S 0.7; 2-phenanthrene m. 135-7°, S 0.5; 3-isomer m. 133-5°, S 0.5; 1-pyrrole m. 85-7°, S 0.9; 2-thiophene m. 66-7°, S 1.8; 2-furan oil, S 0.4; 2,6-dihydroxy-5-pyrimidine m. 271-2°, S 1; 2-methyl-4-hydroxy-5-pyrimidine m. 184°, S 0.9; 3,4-methylenedioxypyrenyl m. 99-100°, S 1; 2-methyl-4-thiazole m.

10/513699

93-4°, S 0.85; 4-methyl-2-thiazole m. 80-2°, S 0.9;
2-pyridine m. 93-4°, S 1; 3-isomer m. 94° S 1;
6-methyl-2-pyridine m. 49-50°, S 1; 2-benzyl-1-imidazole m.
177-9°, S 1; 3-quinoline m. 151-2°, S 1; 6-isomer m.
135°, S 1; 8-isomer m. 92-3°, S 1; 2-benzimidazole m.
185-90°, S 1; 5-isomer m. 160-2°, S 1;
2-hydroxy-5-benzimidazole m. 245-6°, S 1; 7-hydroxy-4-coumarin m.
114-16°, S 1; 9-xanthene m. 157-8°, S 0.8; 9-thiaxanthene m.
148-9°, S 0.7; 5-hydantoin m. 160-2°, S 0.9. Only a few of
these compds. appeared to be utilized readily by the mold for the
formation of new penicillins. Several of the compds. appeared to effect
some increase in penicillin yield or to change the differential assay
value of the crude penicillin produced in their presence.

IT 858199-36-7P, 2-Naphthaleneacetamide,
5,6,7,8-tetrahydro-N-2-hydroxyethyl-
RL: PREP (Preparation)
(preparation of)
RN 858199-36-7 CAPLUS
CN 2-Naphthaleneacetamide, 5,6,7,8-tetrahydro-N-(2-hydroxyethyl)- (CA INDEX
NAME)



OS.CITING REF COUNT: 4 THERE ARE 4 CAPLUS RECORDS THAT CITE THIS RECORD
(4 CITINGS)

L5 ANSWER 34 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1949:11510 CAPLUS

DOCUMENT NUMBER: 43:11510

ORIGINAL REFERENCE NO.: 43:2319i,2320a-c

TITLE: Adrenergic blocking drugs. II. Antagonism of histamine and adrenaline with N-(2-haloalkyl)-1-naphthylmethylanine derivatives

AUTHOR(S): Loew, Earl R.; Micetich, Audrey

SOURCE: Journal of Pharmacology and Experimental Therapeutics (1948), 94, 339-49
CODEN: JPETAB; ISSN: 0022-3565

DOCUMENT TYPE: Journal

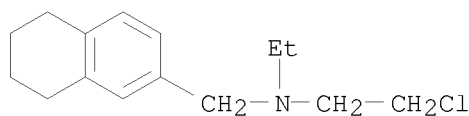
LANGUAGE: Unavailable

AB cf. C.A. 43, 311a. Compds. studied were the series 1-C10H7CH2N(R)CH2CH2Cl where R = Me, Et, Pr, iso-Pr, allyl, Bu, sec-Bu, iso-Bu, Am, hexyl, 2-methoxyethyl, and 2-chloroethyl, and also N-ethyl-N-2-hydroxyethyl-1-naphthylmethylanine, N-ethyl-N-2-chloroethyl-1-(4-chloronaphthyl)methylanine, and N-ethyl-N-2-chloroethyl-2-naphthylmethylanine and its 5,6,7,8-tetrahydro derivative; all as HCl salts. They exhibited the dual property of strongly blocking certain effects of both adrenaline and its physiol. antagonist histamine. The lower alkyl homologs in oral doses of 3-17 mg./kg. were effective in reducing the toxicity of adrenaline in mice; the toxic doses were 60-360 times as great. Injected i.v. in dogs, the compds. reversed the action of adrenaline and diminished the pressor response to injected histamine; their effect was of long duration. S.c. in guinea pigs, they reduced the toxicity of histamine aerosol and the histamine released during anaphylaxis. The most effective compds. were N-ethyl-N-2-chloroethyl- and N-ethyl-N-2-bromoethyl-1-naphthylmethylanine (SY-14 and SY-28).

IT 856200-34-5, 2-Naphthalenemethylanine,
N-(2-chloroethyl)-N-ethyl-5,6,7,8-tetrahydro-, hydrochloride
(antagonism to adrenaline and histamine)

RN 856200-34-5 CAPLUS

CN 2-Naphthalenemethanamine, N-(2-chloroethyl)-N-ethyl-5,6,7,8-tetrahydro-, hydrochloride (1:1) (CA INDEX NAME)



● HCl

10/513699

L5 ANSWER 35 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1948:19976 CAPLUS

DOCUMENT NUMBER: 42:19976

ORIGINAL REFERENCE NO.: 42:4302i,4303a

TITLE: Thermally vaporizable fumigant comprising sensitized ammonium nitrate and a pesticide

INVENTOR(S): Flanders, John Stocks; Jones, Elwyn

PATENT ASSIGNEE(S): Imperial Chemical Industries Ltd.

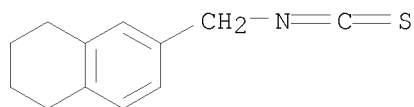
DOCUMENT TYPE: Patent

LANGUAGE: Unavailable

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
	-----	----	-----	-----	-----
	US 2440082		19480420	US 1946-658882	19460401 <--
AB	Ten parts of a pesticide (DDT, γ -hexachlorobenzene, or pentachlorophenol) is incorporated into a mixture of 90 parts of NH_4NO_3 and 10 parts of a chromate, which by its self-sustained exothermic reaction will evaporate the pesticide, after being set off with a fuse.				
IT	855737-77-8P, Naphthalene, 1,2,3,4-tetrahydro-6-(isothiocyanatomethyl)- RL: PREP (Preparation) (preparation of)				
RN	855737-77-8 CAPLUS				
CN	Naphthalene, 1,2,3,4-tetrahydro-6-(isothiocyanatomethyl)- (CA INDEX NAME)				

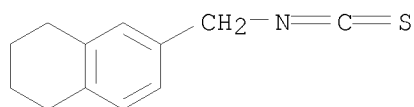


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L5 ANSWER 36 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1948:19975 CAPLUS
DOCUMENT NUMBER: 42:19975
ORIGINAL REFERENCE NO.: 42:4302g-i
TITLE: Tetralyl compounds
INVENTOR(S): Jones, Franklin D.
PATENT ASSIGNEE(S): American Chemical Paint Co.
DOCUMENT TYPE: Patent
LANGUAGE: Unavailable
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
	-----	----	-----	-----	-----
	US 2438751		19480330	US	<--
AB	From tetralyl-6-methyl chloride (I) are prepared tetralyl-6-acetonitrile (also called 1,2,3,4-tetrahydronaphthalene-6-acetonitrile), b. 150-5° at 2 mm.; tetralyl-6-acetic acid, m. 109-14°; the amide, m. 147°; the Me ester, b. 135-40° at 1-2 mm.; the Et ester, b. 140-5° at 1-2 mm.; and alkaline salts, soluble in H2O. On refluxing I and a substantially equivalent mol. proportion of KCNS and 3-4 times their volume of an alc. for 1 hr. and pouring into 4 times its volume of cold H2O, tetralyl-6-methyl thiocyanate is precipitated which, on distillation, is converted to the isothiocyanate, b. 168-74° at 1-2 mm. These derivs. are effective as plant hormones, insecticides, and fungicides.				
IT	855737-77-8P, Naphthalene, 1,2,3,4-tetrahydro-6-(isothiocyanatomethyl)- RL: PREP (Preparation) (preparation of)				
RN	855737-77-8 CAPLUS				
CN	Naphthalene, 1,2,3,4-tetrahydro-6-(isothiocyanatomethyl)- (CA INDEX NAME)				



L5 ANSWER 37 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1946:21658 CAPLUS
 DOCUMENT NUMBER: 40:21658
 ORIGINAL REFERENCE NO.: 40:4234f-i,4235a-d
 TITLE: Aralkylated sulfonamides
 INVENTOR(S): Albrecht, Otto
 PATENT ASSIGNEE(S): Society of Chemical Industry
 DOCUMENT TYPE: Patent
 LANGUAGE: Unavailable
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 2398990		19460423	US 1943-474900	19430205 <--

GI For diagram(s), see printed CA Issue.
 AB Useful wetting, dispersing, and washing agents result from condensation of sulfonamides with aryl chloromethyl compds. Solubility in H2O is achieved by polar groups, such as -SO3Na or -OSO3H. Thus p-cymene is converted with ClSO3H, followed by NH3, to a mixture of cymenesulfonamides, 53 parts of which is treated with 50 parts of Na formaldehyde bisulfite (I) and 2.5 parts of diamylamine for 20 min. at 160-5°; 7 parts of the product is dissolved in 14 parts of H2O and 3.2 parts of 6-(chloromethyl)-1,2,3,4-tetrahydronaphthalene (II) and 2.9 parts of 30% NaOH is added at 65-70° during 1 hr. The mixture is stirred, excess NaOH is neutralized, and by evaporation to dryness a product is isolated, probably of the formula Likewise 10 parts of the product obtained by the reaction of the tetrahydronaphthalenesulfonamides (III) with I is dissolved in 20 parts of H2O and 4.4 parts of II is added and heated with 3.9 parts of 30% NaOH until a sample dissolves in H2O (2 hrs. at 65-70°). Similar products may be obtained with Na formaldehydesulfoxylate, and with a (chloromethyl)cymene (IV), or from the complex mixture which results when tetrahydronaphthalene is heated with AlCl3 at 100°. A technical grade of III will react in aqueous alkali with CH2ClCO2H to yield tetrahydronaphthalene-sulfonamidoacetic acids which react similarly with II to give a soluble washing powder. Another variation consists in preparation of the isomeric N-hydroxyethyl-p-cymenesulfonamides (from the sulfonyl chloride and CH2(OH)CH2NH2) and treating them with IV, followed by treatment with ClSO3H, and then H2O, to give the acid sulfate. Also, 15 parts of the condensation product (V) from the Na salt of 2-amino-6,8-naphthalenedisulfonic acid (VI) and N-chloroacetyltetrahydronaphthalenesulfonamide is dissolved in 25 parts of H2O at 70° and 4.5 parts of II and 3.7 parts of 30% NaOH are added; stirring at 65-70° for 30 min. gives a H2O-soluble product which, however, reacts further with 4.5 parts of II, condensation apparently taking place at both the amido and amino H. V can be made as follows: 46 parts of III is heated with 24.6 parts of CH2ClCOCl to 100° in the course of 2 hrs., and heating is continued for 2.5 hrs. at 100°; 14.4 parts of the chloroacetyl derivative is mixed with 6.7 parts of 30% NaOH and 7.5 parts by volume of EtOH and is dropped during 2 hrs. at 65-70° into a solution of 24.3 parts of VI (partly neutralized as the Na acid salt), containing 62.4% free disulfonic acid, in 50 parts by volume of H2O, made neutral with Na2CO3; V was isolated by evaporation of the reaction mixture to dryness, after it had been stirred for 2.5 hrs. at 70°.
 IT 854747-58-3P, Methanesulfonic acid,
 [N-(5,6,7,8-tetrahydro-2-naphthylmethyl) (5,6,7,8-

10/513699

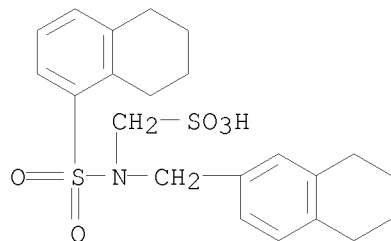
tetrahydronaphthyl)sulfonamido]-, sodium salt 859980-42-0P,
Glycine, N-(5,6,7,8-tetrahydro-2-naphthylmethyl)-N-(5,6,7,8-
tetrahydronaphthylsulfonyl)-, sodium salt

RL: PREP (Preparation)

(preparation of)

RN 854747-58-3 CAPLUS

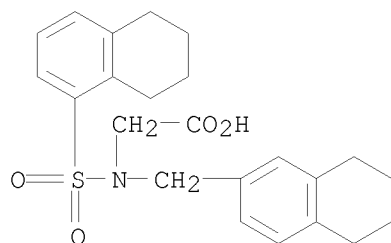
CN Methanesulfonic acid, 1-[[(5,6,7,8-tetrahydro-2-
naphthalenyl)methyl][(5,6,7,8-tetrahydro-1-naphthalenyl)sulfonyl]amino]-,
sodium salt (1:1) (CA INDEX NAME)



● Na

RN 859980-42-0 CAPLUS

CN Glycine, N-(5,6,7,8-tetrahydro-2-naphthylmethyl)-N-(5,6,7,8-
tetrahydronaphthylsulfonyl)-, sodium salt (4CI) (CA INDEX NAME)



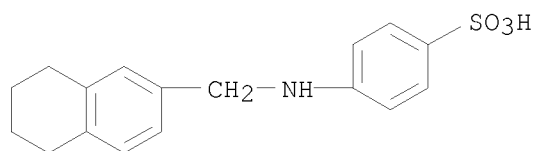
● Na

L5 ANSWER 38 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1946:4643 CAPLUS
 DOCUMENT NUMBER: 40:4643
 ORIGINAL REFERENCE NO.: 40:754i,755a-c,756a-b
 TITLE: Derivatives of sulfonated amines
 INVENTOR(S): Granacher, Charles; Streuli, Paul; Meyer, Jules
 PATENT ASSIGNEE(S): Soc. pour l'ind. chim. a Bale
 DOCUMENT TYPE: Patent
 LANGUAGE: Unavailable
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
	-----	----	-----	-----	-----
	US 2376911		19450529	US 1942-462988	19421022 <--
AB	<p>Aminosulfonic acids are treated with aralkylating agents containing at least 8 C atoms to form N-aralkylaminosulfonic acids which are useful as wetting, dispersing, washing, softening, leveling, or foaming agents in the treatment of textiles. For example the Na salt of sulfanilic acid in aqueous Na₂CO₃ solution is treated with ar-2-(chloromethyl)tetrahydronaphthalene (I) at 70-80° to form the Na salt of N-(ar-2-tetrahydronaphthylmethyl)sulfanilic acid, which is salted out and dried. Similar compds. or their salts are prepared in a like manner by treating salts of the following aminosulfonic acids with I: 1,6-naphthylaminesulfonic acid, 2,6,8(and 1,3,6)-naphthylaminedisulfonic acid, and β-aminoethanesulfonic acid. One or both free H atoms in the NH₂ may be replaced in accordance with the amount of aralkylating agent used. Products containing the N-tetrahydronaphthylmethyl radical have excellent washing properties, and those containing 2 SO₃H groups are particularly good in hard water. Similarly N-(cymylmethyl)aminosulfonic acids and their salts are prepared in a like manner by treating salts of the following aminosulfonic acids with 2-(chloromethyl)cymene: N-methylsulfanilic acid, 2,6,8-naphthylaminodisulfonic acid, phenylhydrazinesulfonic acid, and β-aminoethanesulfonic acid. Again one or both H atoms on the NH₂ group may be replaced. Products containing the cymylmethyl group have particularly good wetting properties. Products containing a free amino H atom may be treated with ethylene oxide or an acyl chloride, such as lauric acid chloride, to form a product with -CH₂CH₂OH (or-CH₂CH₂OCH₂CH₂OH) or an acyl group, resp., on the N atom.</p>				
IT	<p>857954-48-4P, Sulfanilic acid, N-(5,6,7,8-tetrahydro-2-naphthylmethyl)-, sodium salt 857956-72-0P, Taurine, N-(5,6,7,8-tetrahydro-2-naphthylmethyl)-, sodium salt 861090-87-1P, 2-Naphthalenesulfonic acid, 5-(5,6,7,8-tetrahydro-2-naphthylmethylamino)-, sodium salt RL: PREP (Preparation) (preparation of)</p>				
RN	857954-48-4 CAPLUS				
CN	<p>Benzenesulfonic acid, 4-[[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]amino]-, sodium salt (1:1) (CA INDEX NAME)</p>				

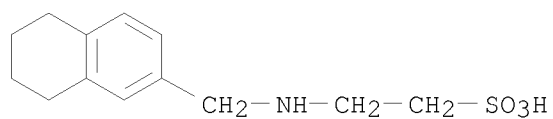
10/513699



● Na

RN 857956-72-0 CAPLUS

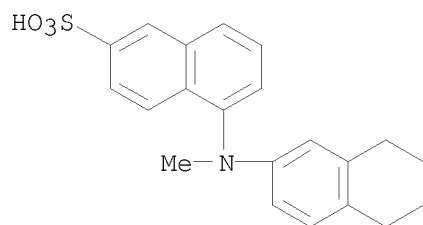
CN Ethanesulfonic acid, 2-[[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]amino]-, sodium salt (1:1) (CA INDEX NAME)



● Na

RN 861090-87-1 CAPLUS

CN 2-Naphthalenesulfonic acid, 5-[methyl(5,6,7,8-tetrahydro-2-naphthalenyl)amino]-, sodium salt (1:1) (CA INDEX NAME)



● Na

L5 ANSWER 39 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1937:53452 CAPLUS

DOCUMENT NUMBER: 31:53452

ORIGINAL REFERENCE NO.: 31:7432i,7433a-i,7434a

TITLE: Friedel-Crafts reaction. I. Synthesis of new compounds in the field of pharmaceutical chemistry

AUTHOR(S): Kranzlein, Paul

SOURCE: Berichte der Deutschen Chemischen Gesellschaft [Abteilung] B: Abhandlungen (1937), 70B, 1776-87

CODEN: BDCBAD; ISSN: 0365-9488

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

OTHER SOURCE(S): CASREACT 31:53452

AB According to Kuhn and his school, in the pigment component of the lactoflavin in vitamin B2 2 o-Me groups in a certain position are responsible for the physiol. action. Flavins not methylated on the benzene nucleus and those in which the Me has been shifted from the 6- to the 5- or from the 7- to the 8-position have no growth effect (C. A. 31, 6239.6). The object of the present work was to synthesize heterocyclic substances containing o-Me groups in corresponding positions. Kunckell and Schneider had observed (C. A. 7, 777) that in the action of ClCH₂COCl on 3,4-Me₂C₆H₃NHAc (I) in the Friedel-Crafts reaction the ClCH₂CO group enters the o-position to the NHAc group, the latter having an o-, the 3-Me group a p-directing influence. It was to be expected that 2-acetamino-5,6,7,8-tetrahydronaphthalene (II) and 5-acetaminohydrindene (III) would behave in the same way, the cyclic methylene residues having the same influence as the 2 o-Me groups in I. Such proved to be the case. 5,6,5',6'-Tetramethylindigo (IV), from 3,4,6-Me₂(ClCH₂CO)C₆H₂NHAc and alkali, was oxidized to 5,6-dimethylisatin (V), best with HNO₃CrO₃ (Ger. pat. 229,815, C. A. 5, 2732). The corresponding dyes were likewise obtained in good yields from the ClCH₂CO derivs. of II and III. The 3 isatins with PhCOMe in alkaline solution gave the 6,7-substituted 2-phenylquinoline-4-carboxylic acids. These acids, as compared with atophan, showed no greater pharmacol. action and about the same, or perhaps somewhat higher, toxicity; they have no vitamin B2 action and have no advantages over other atophan derivs. in their influence on uric acid metabolism. An attempt was also made to introduce the above substituents into acridines. 2'-Chloro-4,5-dimethyl-2-aminobenzophenone (VI) was prepared from I and o-ClC₆H₄COCl but attempts to effect ring closure to the acridone with Cu(OAc)₂ in AmOH and even by heating in PhNO₂ with Cu and K₂CO₃ failed. 3',4'-Dimethyldiphenylamine-2-carboxylic acid (VII) was accordingly prepared by heating 3,4-Me₂C₆H₃NH₂ and o-ClC₆H₄CO₂H with Cu and K₂CO₃; ring closure to 2,3-dimethylacridone (VIII) was easily effected with concentrated H₂SO₄ at 80°. VIII was quant. reduced with Na and AmOH to the dihydroacridine (IX) which with FeCl₃ yielded 2,3-dimethylacridine (X) through a green addition product, IX.X. 2,3-Cyclotrimethylene- (XI) and 2,3-cyclotetramethyleneacridine (XII) were prepared in the same way. The tolerated doses, s.c., per 20-g. mouse weight of X, XI, XII and acridine (XIII) are resp. 40, 20-40, 20 and 2 mg. The dilns. (1:x) at which they inhibit growth in vitro of streptococci are 100, 200, 200, 8000; of staphylococci 200, 200, 200, 4000; of pneumococci 500, <500, 500, 8000. Local disinfection expts. on animals gave similar results; only with gonococci did the new acridines prove nearly as effective as XIII itself. Hence, 2,3-substitution of XIII decreases the toxicity but also the disinfecting power. I (82% from Me₂C₆H₃NH₂ and AcCl in pyridine), m. 96.5°. 6-ClCH₂CO derivative (94%), m. 167°. IV (55%) forms an

olive-yellow Na₂S₂O₄ vat, dissolves in concentrated H₂SO₄ with yellow-red color.

V (72%), orange, m. 214-15°.

2-Phenyl-6,7-dimethylquinoline-4-carboxylic acid (dimethylatophan) (85%), m. 251.5°. II, from the amine and 1.5 mols. Ac₂O on the water bath (80% yield), m. 106°. 3-Chloroacetyl derivative (27%), m. 148°, soluble in concentrated H₂SO₄ with yellow color, developing a strong green fluorescence on short warming. 5,6,5',6'-Bis(cyclotetramethylene)indigo (86%), dark blue. 5,6-Cyclotetramethyleneisatin (85%), brown-orange, m. 194°. 2-Ph - 6,7 - cyclotetramethylenequinoline - 4 - carboxylic acid (53%), m. 237°. III (81%), m. 104°. 6-Chloroacetyl derivative (52%), m. 167°. 5,6,5',6'-Bis(cyclotrimethylene)indigo (83%), soluble in concentrated H₂SO₄ with red color.

5,6-Cyclotrimethyleneisatin

(50%), brown-orange, m. 206°.

2-Phenyl-6,7-cyclotrimethylenequinoline-4-carboxylic acid (40%), light yellow, m. 261°. VI (80%), m. 173°, soluble in concentrated H₂SO₄ with yellow color. VII (83%), m. 188-9°. VIII (80%), yellow, m. 297°, soluble in alc. KOH. IX, m. 215°. X, light yellow, m. 162°, shows green fluorescence in concentrated H₂SO₄.

3',4'-Cyclotetramethylene analog of VII (90%), m. 173°; acridone (78%), yellow, m. 309°; dihydroacridine, yellow, m. 169-70°;

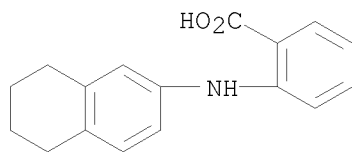
XII, light yellow, m. 117°, shows green fluorescence in H₂SO₄.

3',4'-Cyclotrimethylene analog of VII, m. 176°; acridone (83%), m. 338°, soluble in H₂SO₄ with blue, in MeOH with blue-violet fluorescence; dihydroacridine, m. 209°; XI, m. 152°, shows green fluorescence in H₂SO₄.

IT 856356-54-2P, Anthranilic acid,
N-(5,6,7,8-tetrahydro-2-naphthyl)-
RL: PREP (Preparation)
(preparation of)

RN 856356-54-2 CAPLUS

CN Benzoic acid, 2-[(5,6,7,8-tetrahydro-2-naphthalenyl)amino]- (CA INDEX NAME)



OS.CITING REF COUNT: 2 THERE ARE 2 CAPLUS RECORDS THAT CITE THIS RECORD (2 CITINGS)

L5 ANSWER 40 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1935:22787 CAPLUS

DOCUMENT NUMBER: 29:22787

ORIGINAL REFERENCE NO.: 29:2930i,2931a-e

TITLE: Reduction of nitro and polynitro compounds. XIV. The reduction of aromatic mono-and polynitro compounds

AUTHOR(S): Brand, K.; Mahr, Joseph

SOURCE: Journal fuer Praktische Chemie (Leipzig) (1935

), 142, 153-76

CODEN: JPCEAO; ISSN: 0021-8383

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

AB cf. C. A. 26, 3781. Data are given for the velocity of the reactions
 $2,5\text{-Me}_2\text{C}_6\text{H}_3\text{NO} + 2,5\text{-Me}_2\text{CH}_3\text{N}_6\text{HOH} \rightarrow \text{H}_2\text{O} +$
 $2,5\text{-Me}_2\text{C}_6\text{H}_3\text{N}(:\text{O}):\text{NC}_6\text{H}_3\text{Me}_2\text{-}2,5$; $2,4,5\text{-Me}_2(\text{O}_2\text{N})\text{C}_6\text{H}_2\text{NO} +$
 $2,4,5\text{-Me}_2(\text{O}_2\text{N})\text{C}_6\text{H}_2\text{NHOH} \rightarrow \text{H}_2\text{O} + 2,4,5\text{-Me}_2(\text{O}_2\text{N})\text{C}_6\text{H}_2\text{N}(:\text{O}):$
 $\text{NC}_6\text{H}_2(\text{NO}_2)\text{Me}_2\text{-}2,4,5$; $2,5,3\text{-Me}_2(\text{O}_2\text{N})\text{C}_6\text{H}_2\text{NO} + 2,5,3\text{-Me}_2(\text{O}_2\text{N})\text{C}_6\text{H}_2\text{NHOH}$
 $\rightarrow \text{H}_2\text{O} + 2,5,3\text{-Me}_2(\text{O}_2\text{N})\text{C}_6\text{H}_2\text{N}(:\text{O}):\text{NC}_6\text{H}_2(\text{NO}_2)\text{Me}_2\text{-}2,5,3$; $1\text{-C}_{10}\text{H}_7\text{NO} +$
 $1\text{-C}_{10}\text{H}_7\text{NHOH} \rightarrow \text{H}_2\text{O} + 1,1'\text{-C}_{10}\text{H}_7\text{N}(:\text{O}):\text{NC}_{10}\text{H}_7$; $2\text{-C}_{10}\text{H}_{11}\text{NO} +$
 $2\text{-C}_{10}\text{H}_{11}\text{NHOH} \rightarrow \text{H}_2\text{O} + 2,2'\text{-azoxytetralin}$. The velocity of formation
of azoxybenzene (I) is accelerated by a very small concentration of HO ion; the
effect is much greater than with H ion. Under similar conditions 3,3'-
and 4,4'-azoxytoluene are formed more rapidly but the 2,2'-isomer (II)
much more slowly than I. 2,4,2',4'- and
2,5,2',5'-tetramethylazoxybenzene are formed not only more slowly than I
but also more slowly than II. 3,3'-Dinitroazoxy compds. (III) are formed
considerably more rapidly than the parent compds. The velocity of
formation of III is decreased by the presence of an o-Me group but is
raised by a p-Me group. Cl in the o-position to the NO and NHOH groups
decreases the rate of reaction but in the m- and p-positions it
accelerates it. 2,2'-Dinitroazoxybenzene is formed half as fast, the
3,3'-isomer 5/3 as fast and the 4,4'-isomer 6.5 times as fast as I.
1,1'-Azoxynaphthalene is formed in 70% alc. at a rate only slightly less
than that of I under similar conditions; consts. could not be determined for
acid and alkaline solns., probably because of side reactions. The following
new compds. are reported: 1,2-dimethyl-2-nitro-6-hydroxylaminobenzene,
yellow, m. 87°; 1,4-dimethyl-2-nitro-6-nitrosobenzene, m.
134-5°; 2,5,2',5'-tetramethyl-3,3'-dinitroazoxybenzene, m.
191-2°; 1,3-dimethyl-4-nitro-6-hydroxyl-aminobenzene, yellow, m.
126.5-7.5°; the 6-NO derivative m. 108°; the azoxy compound m.
201-2°; 2-hydroxylamino-tetralin, m. 66-7°;
5,6,7,8,5',6',7',8'-octahydro-2,2'-azoxynaphthalene, yellow, m.
100-1°; the 2,2'-azo derivative, orange-red, m. 127-8°; the
2,2'-hydrazo derivative, pale yellow, m. 121-2°. The theoretical part
discusses many reactions and gives velocity consts. for the formation of
several azoxy compds.

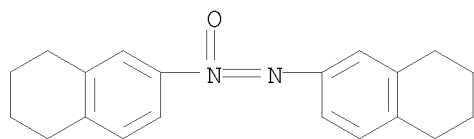
IT 856203-98-0P, Naphthalene, 2,2'-azoxybis[5,6,7,8-tetrahydro-
858024-54-1P, Naphthalene, 2,2'-hydroazobis[5,6,7,8-tetrahydro-
858025-08-8P, Naphthalene, 2,2'-azobis[5,6,7,8-tetrahydro-
RL: PREP (Preparation)

(preparation of)

RN 856203-98-0 CAPLUS

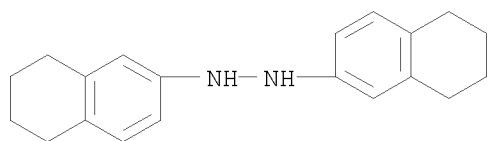
CN Naphthalene, 2,2'-azoxybis[5,6,7,8-tetrahydro- (3CI) (CA INDEX NAME)

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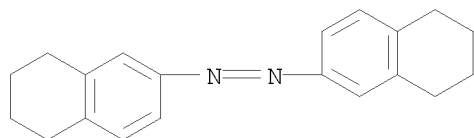
RN 858024-54-1 CAPLUS

CN Naphthalene, 2,2'-hydroazobis[5,6,7,8-tetrahydro- (3CI) (CA INDEX NAME)



RN 858025-08-8 CAPLUS

CN Naphthalene, 2,2'-azobis[5,6,7,8-tetrahydro- (3CI) (CA INDEX NAME)



OS.CITING REF COUNT: 2

THERE ARE 2 CAPLUS RECORDS THAT CITE THIS RECORD
(2 CITINGS)

L5 ANSWER 41 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1935:14013 CAPLUS

DOCUMENT NUMBER: 29:14013

ORIGINAL REFERENCE NO.: 29:1812a-i

TITLE: Homologs of naphthacene. II. 2-Methyl- and 2,7-dimethylnaphthacene; synthetic applications of 2,6-and 2,7-dimethyl-1,2,3,4-tetrahydronaphthalene

AUTHOR(S): Coulson, Edward A.

SOURCE: Journal of the Chemical Society (1935) 77-83

CODEN: JCSOA9; ISSN: 0368-1769

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

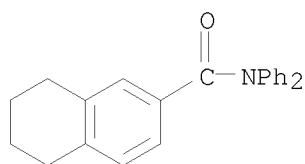
AB cf. C. A. 29, 154.5. 3,6,2-Me₂C₁₀H₅OH (38 g.), 30 g. (NH₄)₂SO₃ and 400 cc. NH₄OH (d. 0.88), heated (6 hrs. at 200°, give 65% of 3,6-dimethyl-2-naphthylamine (I), m. 139°; HCl salt, m. 283° (decomposition); Ac derivative, m. 207°. I (16.5 g.), through the Sandmeyer reaction, gives 11 g. of 3,6-dimethyl-2-naphthonitrile, in. 145°; heating with 50% KOH and EtOH for 24 hrs. gives 11 g. 3,6-dimethyl-2-naphthoic acid, pale cream, m. 224°; acid chloride, cream, m. 70°; the anilide, pale straw, m. 207-8°; the chloride with C₆H₆ or PhMe and AlCl₃, gives resinous products; PhMe and FeCl₃ at 90° for 4 hrs. give a small yield of 2-p-toluyyl-3,6-dimethylnaphthalene, m. 112°; this chars at 400° but forms a small quantity of 2,7-dimethylnaphthacene, golden orange, m. 362°; the solns. show a marked green fluorescence; the cold concentrated H₂SO₄ solution is moss-green. 2,7-Dimethylnaphthacene-9,10-quinone, yellow, m. 223°; the deep purple-red solution in concentrated H₂SO₄ fades on dilution. Diphenylcarbonyl chloride (II), tetralin, AlCl₃ and CS₂, refluxed 3 hrs., give 1,2,3,4-tetrahydro-6-naphthodiphenylamide, m. 87-8°; hydrolysis gives 1,2,3,4-tetrahydro-6-naphthoic acid, m. 154 (acid chloride (III), b₁₂ 163°); 1,2,3,4-tetrahydro-6-naphthanilide, m. 147°. III (28 g.), 30 g. m-C₆H₄Me₂, CS₂ and AlCl₃, refluxed 3 hrs., give 36.5 g. of 6-(2',4'-dimethylbenzoyl)-1,2,3,4-tetrahydronaphthalene (IV), pale yellow, b₁₀ 223°; IV also results in 35.5 g. yield from 30 g. 2,4-Me₂C₆H₃COCl, 30 g. tetralin and 30 g. AlCl₃. Pyrolysis of IV gives a mixture of 2-methylnaphthacene (V), golden orange, m. 350°, 7-methyl-1,2,3,4-tetrahydronaphthacene and 7-methyl-1,2-benzanthracene; the last 2 could not be separated but on dehydrogenation with Se yielded a mixture of V and 7-methyl-1,2-benzanthracene, separated by crystallization from AcOH.

2,6-C₁₀H₆Me₂ on catalytic reduction (Mo catalyst) at 390-400° for 6 hrs. gives 25-30% of the 1,2,3,4-tetrahydro derivative (VI), b. 237-9°, m. 14-17°; there also results some 2,6-dimethyldecalin, b. 216-7°; probably other isomers are formed. 2,7-C₁₀H₆Me₂ (250 g.) at 300° for 4 hrs. gives 84 g. of the 1,2,3,4-tetrahydro derivative (VII), b. 237-8°; a 2,7-dimethyldecalin, b. 216-8°, is also formed. VI and II with AlCl₃ in CS₂ give 2,6-dimethyl-1,2,3,4-tetrahydro-7-naphthoic acid, m. 183°, after hydrolysis of the amide; Se gives 2,6,3-Me₂C₁₀H₅CO₂H; VII yields 2,7-dimethyl-1,2,3,4-tetrahydro-6-naphthoic acid, m. 187°. VI, sulfonated, the Na salt treated with PCl₅ and the chloride with NH₄OH, gives 2,6-dimethyl-1,2,3,4-tetrahydronaphthalene-7-sulfonamide, cream, m. 166-7°. Fusion of the Na salt with KOH at 300-40° gives 2,6-dimethyl-1,2,3,4-tetrahydro-7-naphthol, m. 116°. 2,7-Dimethyl-1,2,3,4-tetrahydro-6-sulfonamide, cream, m. 145.5°; the 6-naphthol m. 87°. 7-p-Toluyyl-2,6-dimethyl-1,2,3,4-

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tetrahydronaphthalene, m. 95°; 6-benzoyl-2,7-dimethyl derivative, b4 202°; 6-p-toluy1 analog, b2 199°. Pyrolysis of these ketones gives: 2,6-dimethyl-1,2,3,4-tetrahydronaphthacene, pale yellow, m. 214°; 2-Me derivative, pale yellow, m. 203°; 2,7-di-Me derivative, pale yellow, m. 210°. Dehydrogenation gives the naphthacene compds.; the 2,6-di-Me derivative is less readily dehydrogenated. Both the tetrahydronaphthols have "wetting-out" properties but the 2,6-isomer is much superior.

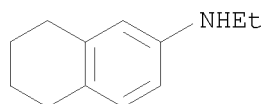
IT 859071-22-0P, 2-Naphthamide, 5,6,7,8-tetrahydro-N,N-diphenyl-
RL: PREP (Preparation)
(preparation of)
RN 859071-22-0 CAPLUS
CN INDEX NAME NOT YET ASSIGNED



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L5 ANSWER 42 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 1928:20295 CAPLUS
DOCUMENT NUMBER: 22:20295
ORIGINAL REFERENCE NO.: 22:2379b-c
TITLE: Hydrogenated naphthylamines
PATENT ASSIGNEE(S): Soc. anon. pour l'ind. chim. a Bale
DOCUMENT TYPE: Patent
LANGUAGE: Unavailable
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

	PATENT NO. -----	KIND ----	DATE -----	APPLICATION NO. -----	DATE -----	
	GB 276571		19270224	GB		<--
AB	ar-Tetrahydronaphthylamine derivs. are made by subjecting N-substituted naphthylamines to catalytic hydrogenation, and the hydrogenated naphthylamines themselves may be made by using an acetyl derivative as the starting material and subsequently saponifying Examples are given for the production of ar-N-ethyltetrahydronaphthylamine, ar-acetyltetrahydro- β -naphthalide, ar-N-phenyltetrahydro- α -naphthylamine and acetylated ar-tetrahydro-N-ethyl- α -naphthylamine.					
IT	856213-39-3P, 2-Naphthylamine, N-ethyl-5,6,7,8-tetrahydro- RL: PREP (Preparation) (preparation of)					
RN	856213-39-3 CAPLUS					
CN	2-Naphthalenamine, N-ethyl-5,6,7,8-tetrahydro- (CA INDEX NAME)					



L5 ANSWER 43 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1924:4947 CAPLUS

DOCUMENT NUMBER: 18:4947

ORIGINAL REFERENCE NO.: 18:675g-i,676a-i

TITLE: Catalytic hydrogenations under pressure in the presence of nickel salts. VI. Nitriles

AUTHOR(S): v. Braun, Julius; Blessing, Georg; Zobel, Friedrich

SOURCE: Berichte der Deutschen Chemischen Gesellschaft

[Abteilung] B: Abhandlungen (1923), 56B, 1988-2001

CODEN: BDCBAD; ISSN: 0365-9488

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

GI For diagram(s), see printed CA Issue.

AB cf. C. A. 17, 2884. With the author's apparatus nitriles can be quickly and smoothly reduced to mixts. of primary and secondary bases (80-95% yields); only in the case of aliphatic nitriles, especially of the lower series, does the reduction soon come to a standstill, probably because the catalyst is paralyzed by slight decomposition products. The pressure has no influence on the nature of the reduction products and the influence of temperature is also not marked and is very variable; to avoid this last factor the present work was almost all done at 115-25°. Two factors whose influence is extraordinarily pronounced are the nature of the solvent and the concentration. The sum total of the reduction products always increases in passing from hydrocarbons like tetra- or decahydronaphthalene to solvents containing O (alcs., ethers), and by varying such O-containing solvents, an extraordinarily marked shifting of the yield in favor of the primary or of the secondary base can be effected. In all solvents, increasing concentration favors the formation of the primary base, often to a very considerable extent. The mechanism of the reaction is probably as follows: $\text{RCN} \rightarrow \text{RCH:NH} \rightarrow \text{RCH}_2\text{NH}_2$; $\text{RCH:NH} + \text{RCH}_2\text{NH}_2$ o-Substitution products of PhCN give, under the same conditions, less secondary base than m- and p-derivs., α -tetralyl cyanide gives less than the β -isomer. When an alc., R'OH, with an especially mobile HO group (PhCH₂OH, cyclohexanol) is used as solvent, mixed bases are also formed: $\text{RCH:NH} + \text{R'OH} \rightarrow \text{RCH(OH)NHR'} \rightarrow \text{RCH:NR'} \rightarrow \text{RCH}_2\text{NHR'}$. The Ni salt was reduced in an autoclave in the desired solvent, then the nitrile, in the amount of solvent necessary to give the desired concentration in the mixed solution, was drawn in and the reduction effected under an average excess pressure of 20 atmospheric. The H was absorbed at the rate of 1 l. in 2-6 min. Heptyl cyanide, b15 87-8°, in tetralin or decalin gives in 25% solution 15 and 18%, in 70% solution 17 and 21%, resp., of octylamine, b14 72-3°, and dioctylamine, b14 175°. PhO(CH₂)₃CN (I) in tetralin (24%) gives 29% PhO(CH₂)₄NH₂, b12 140°, and 47% de- δ -phenoxybutylamine, b15 266°, m. 51-2° (HCl salt, m. 165°; NO derivative, m. 50°; the picrate, Ac and Bz derivs. are oils); heated several hrs. at 100° with fuming HBr, the sec. amine yields di- δ -bromobutylamine dihydrobromide (II), m. 200°, whose aqueous solution, treated with exactly 2 mols. NaOH, almost immediately becomes neutral and clear; on evaporating, extracting with CHCl₃

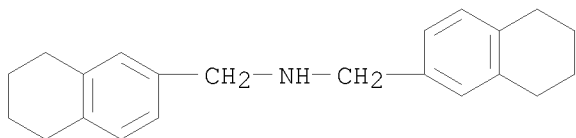
and

adding Et₂O there is at once precipitated the very hygroscopic bispyrrolidinium bromide, C₈H₁₇NBr₃, m. 256-8°; the primary product, N- δ -bromobutylpyrrolidine, (CH₂.CH₂)₂N(CH₂)₄Br, formed by intramol. alkylation of the free base of II, cannot be isolated even when the reaction mixture is carefully cooled. When the I is reduced in cyclohexanol (17-25% solution) there is formed, in addition to 37 and 30%, resp., of the

above primary and sec. bases, 15% of the mixed cyclohexyl- γ -phenoxybutylamine, b16 177-9°, isolated as the picrate, m. 110°; the HCl salt is deliquescent and the NO derivative oily. The yields of PhCH₂NH₂ (III), b13 75-80°, and of (PhCH₂)₂NH (IV), b13 160-5°, resp., from PhCN in various solvents (% concentration of the solution in parentheses) are as follows: decalin or tetralin (9) 44, 40, (25) 41, 35 (66) 72, 5; EtOH (9) 59, 14, (66) 71, 8; Am₂O (20) 41, 39; in cyclohexanol (16%) in addition to 24 and 11% of III and IV is obtained 35% cyclohexylbenzylamine, b15 145-7° (HCl salt, m. 284°; NO derivative, m. 43°; PhSO₂ derivative, m. 90°); in m-methylcyclohexanol (15%) are obtained 58 and 15% of III and IV and 5% m-methylcyclohexylbenzylamine, b15 155° (NO derivative, oil; HCl salt, m. 249°; HBr salt, m. 250°). The C₁₀H₇CN have to be reduced at 190° in order to absorb the H with reasonable rapidity. The α -compound in decalin or tetralin (45%) gives 70% α -naphthylmethylamine, b12 155° (HCl salt, m. 262-4°; picrate, m. 223°; phenylurea, m. 216°; Ac derivative, m. 134°; PhSO₂ derivative, m. 148°; quaternary MeI salt, m. 213°), and 21% of di- α -naphthylmethylamine, m. 73-4°, isolated as the HCl salt, m. 239°; picrate, m. 202°; NO derivative, m. 147°; quaternary MeI derivative, m. 209-10°. β -C₁₀H₇CN in 50% solution gives 66% β -naphthylmethylamine, b12 148-9°, m. 60° (HCl salt, m. 269°; picrate, m. 226°; Ac derivative, m. 126°; quaternary methiodide, m. 168°), and 17% of di- β -naphthylmethylamine, m. 95% (HCl salt, m. 285°; picrate, m. 126°; NO derivative, m. 132°; quaternary methiodide, m. 217°). The yields of PhCH₂CH₂NH₂ (V) and (PhCH₂CH₂)₂NH (VI), b18 195°, m. 28-30° (picrate, m. 150°; NO derivative, m. 53°; phenylthiourea, m. 113°) from PhCH₂CN in various solvents (concentration of solution in parentheses) are as follows: tetralin or decalin (20) 39, 21, (23) 35, 27, (33) 36, 26, (66) 64, 3; EtOH (14) 8, 78, (25) 20, 60, (50) 22, 62; octyl alc. (50) 55, 38; Ph(CH₂)₂OH (50) 55, 35; ac- β -tetralol (50) 23, 46; cyclopentanol (20) 71, 17; Am₂O (15) 13, 75, (66) 57, 29; in cyclohexanol (15%) are obtained 35 and 10% V and VI and 38% β -phenylethylcyclohexylamine, b13 163-9° (HCl salt, m. 199°; picrate, m. 154°); in PhCH₂OH (20%), 61% V and 26% β -phenylethylbenzylamine, b15 186-7° (HCl salt, m. 254°; NO derivative, m. 142°; Bz derivative, m. 123°; picrate, m. 146°); in p-MeC₆H₄CH₂OH (20%), 45% V and 28% β -phenylethyl-p-methylbenzylamine, b14 191-3° (HCl salt, m. 238-40°; picrate, m. 139-41°). PhCH₂CH₂CN in decalin or tetralin (33%) gives 57 and 29%, in Ph(CH₂)₂OH (16%) 70 and 20%, resp., of Ph(CH₂)₃NH₂ (VII), b18 112-4°, and (PhCH₂CH₂CH₂)₂NH, b18 220-2°; in PhCH₂OH (10%) 15% VII and 45% β -phenylpropylbenzylamine, isolated as the HCl salt m. 184-5°. o-MeC₆H₄CN gives in decalin or tetralin (10) 54, 32, (81) 80, 9, in EtOH (14) 72 and 16%, resp., of MeC₆H₄CH₂NH₂ and di-o-methylbenzylamine, b16 190° (HCl salt, m. 202°; picrate, m. 133°). m-MeC₆H₄CN in decalin or tetralin (19, 37 and 82%) yields 54, 70 and 75%, resp., of MeC₆H₄CH₂NH₂ and 32, 14 and 15% of di-m-methylbenzylamine, b14 189-91° (HCl salt, m. 199°; Bz derivative, m. 100°). From p-MeC₆H₄CN (30% solution) are obtained 41% MeC₆H₄CH₂NH₂ and 32% di-p-methylbenzylamine, b30 220°. o-MeOCH₂C₆H₄CN in 20% solution gives 44 and 22, the p-compound in 50% solution 20 and 24%, resp., of the primary and secondary bases (cf. C. A. 17, 2582). α -Tetralyl cyanide in 20% solution gives 70 and 1.5%, resp., of ar- α -tetralylmethylamine, b14

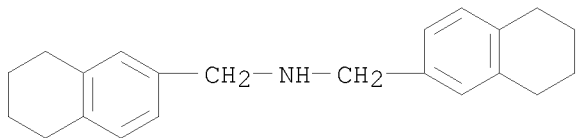
150°, and of the sec. base, m. 93° (HCl salt, m. 212°; NO derivative, m. 90-1°), while the β -isomer in 30% solution yields 47% of the primary base, b11 147°, and 24% of the sec. base, b11 265-7° (HCl salt, m. 245°; Bz derivative, m. 241-2°; NO derivative, m. 76°).

IT 861318-28-7P, Dimethylamine,
 α, α' -bis(5,6,7,8-tetrahydro-2-naphthyl)-, -HCl
 861375-66-8P, Dimethylamine,
 α, α' -bis(5,6,7,8-tetrahydro-2-naphthyl)-
 861376-22-9P, Dimethylamine,
 N-nitroso- α, α' -bis(5,6,7,8-tetrahydro-2-naphthyl)-
 861787-07-7P, Benzamide, N,N-bis(5,6,7,8-tetrahydro-2-naphthylmethyl)-
 RL: PREP (Preparation)
 (preparation of)
 RN 861318-28-7 CAPLUS
 CN 2-Naphthalenemethanamine, 5,6,7,8-tetrahydro-N-[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]-, hydrochloride (1:1) (CA INDEX NAME)

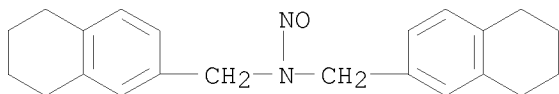


● HCl

RN 861375-66-8 CAPLUS
 CN 2-Naphthalenemethanamine, 5,6,7,8-tetrahydro-N-[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]- (CA INDEX NAME)

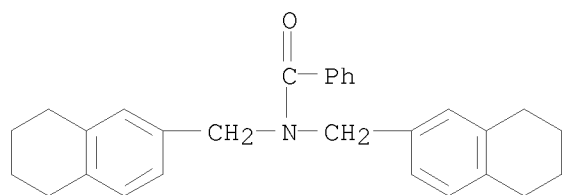


RN 861376-22-9 CAPLUS
 CN 2-Naphthalenemethanamine, 5,6,7,8-tetrahydro-N-nitroso-N-[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]- (CA INDEX NAME)



RN 861787-07-7 CAPLUS
 CN Benzamide, N,N-bis[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]- (CA INDEX NAME)

10/513699



<12/04/2007>

Erich Leese

L5 ANSWER 44 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1922:24682 CAPLUS

DOCUMENT NUMBER: 16:24682

ORIGINAL REFERENCE NO.: 16:4202d-i, 4203a

TITLE: Benzopolymethylene compounds. IV. The two ar-aldehydes of tetralin

AUTHOR(S): Braun, Julius V.; Moldaenke, K.; Dirlam, H.; Gruber, H.

SOURCE: Berichte der Deutschen Chemischen Gesellschaft [Abteilung] B: Abhandlungen (1922), 55B, 1700-9

CODEN: BDCBAD; ISSN: 0365-9488

DOCUMENT TYPE: Journal

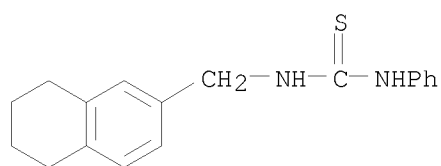
LANGUAGE: Unavailable

AB When tetralin (A) is treated with CO and HCl in the presence of AlCl₃ it is impossible to prevent the greater part of the A from condensing with itself in the same way as it does with AlCl₃ alone (Schroeter, C. A. 15, 525); the small part that does react with the CO and HCl gives exclusively the ar-tetralin- β -aldehyde (B) (2 g. from 100 g. A). Both B and the α -isomer (C) can be obtained from the ar-tetralin- β - and α -methylamines (D and E, resp.) through the corresponding alcs. Bamberger and Lodter's statement that α -C₁₀H₇CH₂NH₂ on reduction takes up the H in the substituted nucleus (Ber. 20, 1708(1887)) seemed to exclude this compound and the β -isomer as the starting points in the synthesis of B and C. Accordingly the NH₂ group in the α - and β -tetralylamines was replaced by CN, which was then reduced to CH₂NH₂, but the yields are poor. On repeating B.'s work, however, it was found that it is the unsubstituted nucleus which takes up the H on reduction and that the pure D and E can easily be obtained in this way. ar- α -Tetralyl cyanide, obtained in 22% yield from the amine by the Sandmeyer reaction, b₁₅ 153°, solidifies to a yellowish crystalline mass m. 48° (Bamberger and Bordt, Ber. 22, 625(1889), describe it as an oil b₁₂₁ 277-9°, which does not solidify), hydrolyzed by fuming HCl in a sealed tube at 120° to the acid, m. 150° (B. and H. give 123°); reduction of the nitrile with Na and alc. gives chiefly A and only about 1/3 is converted into E, oil of basic odor, b₁₁ 149-52°, eagerly absorbs CO₂ from the air, also obtained in almost 90% yield from α -C₁₀H₇CH₂NH₂ with 8 atoms of Na in AmOH (in EtOH there is very little reaction); hydrochloride, silvery needles from alc., m. 253°; picrate, golden yellow prisms from alc., m. 242°; acetyl derivative, m. 125°; benzoyl derivative, m. 144°; phenylurea, m. 199°; phenylthiourea, m. 153°. The corresponding ar- α -tetralylmethylamine (from α -C₁₀H₇CN with Na and alc.) forms a hydrochloride m. 230°, picrate m. 169-70°, phenylurea m. 126°, and benzoyl derivative m. 125°. ar- β -Tetralyl cyanide (obtained in 45-60% yield), liquid of a not unpleasant odor, b₁₁ 151-2°, m. 20-1°, gives with Na and EtOH 30% of D, b₁₁ 146-8°; hydrochloride, m. 248°; picrate, m. 215°; benzoyl derivative, long needles from alc., m. 165°, b₁₀ 260-5°; p-nitrobenzoyl derivative, m. 170°; phenylthiourea, m. 130°. D is also obtained in almost 90% yield from β -C₁₀H₇CH₂NH₂ with Na and AmOH. ar- α -Tetralylcarbinol, obtained in 80% yield from E diazotized in AcOH with the calculated amount of NaNO₂ and heated on the H₂O bath until the evolution of gas ceases, b₁₂ 154-5°, gives in H₂SO₄ with the calculated amount of K₂Cr₂O₇ 1/3 of its weight of C, b₁₂ 131-3°, as an almost odorless oil; semicarbazone, m. 187°. KMnO₄ smoothly oxidizes C to the acid.

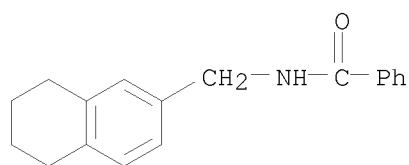
10/513699

ar- β -Tetralylcarbinol (yield, 70%), faintly yellow liquid with a strong pleasant odor, b₁₄ 148-52°, gives on oxidation 25% of B, liquid of characteristic peppermint-like odor, b₁₄ 138°; semicarbazone, m. 219°.

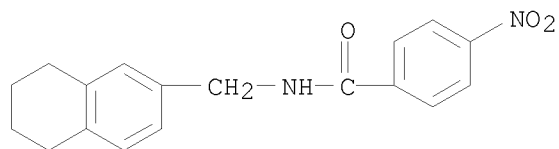
IT 861521-58-6P, Urea, α -phenyl- β -[(5,6,7,8-tetrahydro-2-naphthyl)methyl]thio- 861800-58-0P, Benzamide, N-[(5,6,7,8-tetrahydro-2-naphthyl)methyl]- 861969-08-6P, Benzamide, p-nitro-N-[(5,6,7,8-tetrahydro-2-naphthyl)methyl]-
RL: PREP (Preparation)
(preparation of)
RN 861521-58-6 CAPLUS
CN Thiourea, N-phenyl-N'-[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]- (CA INDEX NAME)



RN 861800-58-0 CAPLUS
CN Benzamide, N-[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]- (CA INDEX NAME)



RN 861969-08-6 CAPLUS
CN Benzamide, 4-nitro-N-[(5,6,7,8-tetrahydro-2-naphthalenyl)methyl]- (CA INDEX NAME)



L5 ANSWER 45 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1922:10116 CAPLUS

DOCUMENT NUMBER: 16:10116

ORIGINAL REFERENCE NO.: 16:1763h-i,1764a-i,1765a-c

TITLE: Hydrogenated naphthalenes and their transformations.
II. Nitro and amino derivatives of
tetrahydronaphthaleneAUTHOR(S): Schroeter, G.; Kindermann, E.; Dietrich, C.;
Beyschlag, C.; Fleischhauer, Cl.; Riebenschahm, E.;
Oesterlin, C.SOURCE: Justus Liebigs Annalen der Chemie (1922),
426, 17-83

CODEN: JLACBF; ISSN: 0075-4617

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

AB The work described covers the nitration of C₁₀H₁₂, the reduction of various mono-, di- and tri-NO₂ derivs., and the nitration of the Ac derivs. of the amines so obtained. The orientations of a considerable no. of isomeric compds. are definitely established. The mononitration of C₁₀H₁₂, using a mixture of HNO₃ and H₂SO₄, leads to the formation of both 1- and 2-nitro-ar-tetrahydronaphthalene, which may be separated by fractional distillation and "freezing out" the fractions or by taking advantage of the

fact

that the 2-NO₂ compound is more easily reduced than its isomer to an NH₂ derivative 1-Nitroderiv. m. 34°, b₁₃ 157°, d₄₀₄₀ 1.1757, and the 2-NO₂ deriv. m. 31.4°, b₁₃ 169°, d₄₀₄₀ 1.1762. On dinitration, C₁₀H₁₂ yields a mixture of 1,2- and 1,3-dinitro-ar-tetrahydronaphthalene (1,2-derivative, m. 102-3°; 1,3-derivative, m. 95°) which may be separated by crystallization from

concentrated H₂SO₄

in which the former is less soluble. The orientation of the 1,2-compound rests on its reduction (see below) and that of the 1,3-derivative on its oxidation to 3,5-(O₂N)₂C₆H₂(CO₂H)₂ and its reduction. Another oxidation product with HNO₃ is β-o-carboxytrinitrophenylpropionic acid, which decomps. violently on heating. The potassium hydrogen salt was analyzed. The 1,3-derivative cannot be further nitrated. The 1,2-derivative yields 1,2,4-trinitro-ar-tetrahydronaphthalene, m. 94.5-5°, the structure of which was established by reduction. 1,1-Hydrazo-ar-tetrahydronaphthalene, by reduction of the 1-NO₂ derivative with Zn dust and alkali, slender needles, m. 181-3°, and on oxidation with KMnO₄ is converted quant. into 1,1-azo-ar-tetrahydro-naphthalene, glistening red needles, m. 190-1°, also obtained, with the 1,1-azoxyderivative, yellow needles, m. 184°, by reduction of the NO₂ derivative with Zn and NaOH under less energetic conditions. The benzidine conversion gives rise to 4,4'-diamino-1,1'-di-ar-tetrahydronaphthyl, m. 153-4°. The hydrochloride, hydrobromide, sulfate and phosphate are described. The corresponding diazonium salt gives substantive dyes on coupling with various intermediates. 4,4'-Dihydrazino-1,1'-di-ar-tetrahydronaphthyl is formed by reduction with SnCl₂; 4,4'-diethoxyderivative, needles, m. 173°. A compound, probably 1,1'-diamino-2,2'-di-ar-tetrahydronaphthyl, results as a by-product in the preparation of the 4,4'-derivative, needles, m. 216°; on heating the HCl salt, it yields NH₄Cl and a carbazole-like base, separated as the picrate.

α- and β-C₁₀H₁₁NH₂ are obtained from the corresponding NO₂ compds. by catalytic reduction and may also be obtained by reducing the crude mononitration product of C₁₀H₁₂ and separating the isomeric bases by the differences in solubility of their HCl salts in H₂O, their methanedisulfonates

in 96% EtOH and the difference in the ease with which the bases react with Ac2O. α -C10H11NH2.HCl is more easily soluble in H2O than the β -derivative. The methanedisulfonate, leaflets, is soluble in 20 parts hot H2O, 60 parts cold H2O and 6 times as soluble in EtOH as the β -derivative. C6H4(CO)2O gives α -tetralylphthalamic acid, needles, m. 182-4° (decomposition), which loses H2O on heating and gives the imide, long needles, m. 148-50°. The action of Me2SO4 on the Ac derivative gives α -acetmethylaminotetralin, needles, m. 70-2°, b11 182-5°. β -Aminotetralin hydrochloride, large leaflets, is sparingly soluble in cold H2O, as is the sulfate; the methanedithionate forms leaflets which are sparingly soluble in alc. The phthalamic acid forms glistening needles, m. 156.5-8.5°, and the phthalimide, needles, m. 169-71°. β -Acetmethylaminotetralin, needles, m. 67-9°, b12 178-80°. The nitration of α -C10H11NHAc yields 1,4-C10H10(NH2)NO2 (Green and Rowe, C. A. 13, 710) and as by-product, 1-acetamino-2-nitrotetralin, needles m. 184-5°, and 1-acetamino-3-nitrotetralin, needles, m. 193°. 1-Amino-2-nitrotetralin, by saponification of the Ac derivative, forms long orange needles, m. 87-8°. 3-Nitro derivative, yellow leaflets, m. 78°, is also obtained by regulated reduction of the 1,3-di-NO2 derivative. The nitration of β -C10H11NHAc in AcOH gives as the main product 2-acetamino-3-nitrotetralin, long yellow needles, m. 134-35.5°, while 2-amino-1-nitrotetralin, needles, m. 128-9°, is produced only in small amts. In H2SO4, the main product is 2-acetamino-4-nitrotetralin, long needles, m. 194°, with the 3-NO2 derivative as a by-product. 2-Amino-3-nitrotetralin, long red needles, m. 125-7°. Me2SO4 gives the methyl derivative, fine red needles, m. 113-5°, which, with Ac2O, gives an acetyl derivative, AcNMeC10H10NO2, m. 107-8.5°. 2-Amino-1-nitrotetralin, red needles, m. 96°, obtained by hydrolysis of the Ac derivative and also by partial reduction of the 1,2-(NO2)2 derivative. 2-Amino-4-nitrotetralin, yellow, m. 55°, which may be diazotized and which yields 4-nitro-2-hydroxytetrahydronaphthalene, amorphous body. The diazo compound may be easily reduced to 1-O2NC10H11. 1,3-Dinitro-2-acetaminotetralin is formed by the further nitration of the 3-NO2 derivative, needles, m. 189-91°. 1,3-Dinitro-2-aminotetralin, yellow needles, m. 166-8°. 3,4-Dinitro-2-acetaminotetralin, needles, m. 175-7°, yields, on hydrolysis, 3,4-dinitro-2-aminotetralin, long golden yellow needles, m. 157°. Tetrahydro-2,3-naphthylenediamine, by catalytic reduction of the 3-NO2 derivative with H, glistening leaflets, m. 135-6°, b13 165°. The hydrochloride forms glistening leaflets. With AcOH it forms 2,3-tetralylene- μ -methylimidazole, m. 251-2°, and with phenanthrenquinone 2,3-tetralylenephenanthrazine, pale yellow, glistening needles, m. 214-6°. 1-Acetamino-2-aminotetralin, m. 149-51°, yields with Ac2O the 1,2-diacetate, m. 244-5°; 1,2-tetralylphenanthrazine, small, light yellow needles, m. 230°. 1,3-Diaminotetralin forms pearly leaflets, m. 84-5°, the 3-acetate of which forms glistening needles, m. 110-1°, the 1-acetate, m. 173°, and the diacetate, small needles, m. 245-6°. Monoacetyl-1,4-diaminotetralin, glistening needles, m. 154-6°. 1,2,3-Triaminotetralin, by reduction of the 3,4-(NO2)2 or the 1,3-(NO2)2 derivative, is unstable in air but gives a crystalline hydrochloride, and a triacetate, needles, m. 285°. 1,2,4-Triaminotetralin, also unstable, forms a triacetate, fine needles, m. 315°.

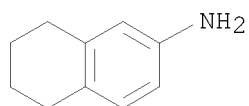
IT 861352-73-0P, 2-Naphthylamine, 5,6,7,8-tetrahydro-, methanedisulfonate

10/513699

RL: PREP (Preparation)
(preparation of)
RN 861352-73-0 CAPLUS
CN Methanedisulfonic acid, compd. with 5,6,7,8-tetrahydro-2-naphthalenamine
(1:1) (CA INDEX NAME)

CM 1

CRN 2217-43-8
CMF C10 H13 N



CM 2

CRN 503-40-2
CMF C H4 O6 S2

HO₃S-CH₂-SO₃H

OS.CITING REF COUNT: 2 THERE ARE 2 CAPLUS RECORDS THAT CITE THIS RECORD
(2 CITINGS)

L5 ANSWER 46 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1919:229 CAPLUS

DOCUMENT NUMBER: 13:229

ORIGINAL REFERENCE NO.: 13:43b-e

TITLE: Transformation of tetrahydronaphthalene (tetralin) in the animal body

AUTHOR(S): Schroeter, G.; Thomas, K.

SOURCE: Journal of the Chemical Society, Abstracts (1918), 114(I), 418

CODEN: JCSAAZ; ISSN: 0590-9791

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

AB Tetrahydronaphthalene fed to a dog is absorbed and eventually excreted in combination with urea. The existence of 4 compds. of tetrahydronaphthalene and urea is theoretically possible, and the authors have succeeded in preparing all of them by the interaction of KCNO on the respective amines in H₂O. ar-Tetrahydro- α -carbamidonaphthalene, C₁₁H₁₄ON₂, crystallizes in square plates from alc., soften at 198° and melts at about 206° (quickly heated, at 212°). ar-Tetrahydro- β -carbamidonaphthalene, needles, m. 134° (decomposition). ac-Tetrahydro- α -carbamidonaphthalene, needles, m. 210.5°. ac-Tetrahydro- β -carbamidonaphthalene, needles, m. 183°. Comparison of the natural product with these compds. shows that tetralin in its passage through the body is converted into dl-ac-tetrahydro- α -carbamidonaphthalene. In the preparation of the ar- β -compds., a small amount of a substance was obtained in the form of needles, which did not melt below 245° and possessed the composition of di-ar-tetrahydro- β -naphthylcarbamide, (C₁₀H₁₂N)₂CO.

IT 871892-48-7P, Urea, s-bis(5,6,7,8-tetrahydro-2-naphthyl)-

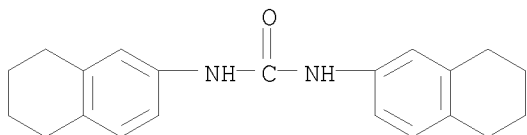
872283-39-1P, Urea, (5,6,7,8-tetrahydro-2-naphthyl)-

RL: PREP (Preparation)

(preparation of)

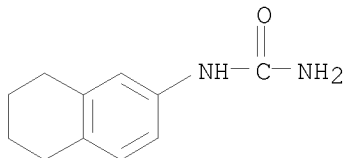
RN 871892-48-7 CAPLUS

CN Urea, N,N'-bis(5,6,7,8-tetrahydro-2-naphthalenyl)- (CA INDEX NAME)



RN 872283-39-1 CAPLUS

CN Urea, N-(5,6,7,8-tetrahydro-2-naphthalenyl)- (CA INDEX NAME)



L5 ANSWER 47 OF 47 CAPLUS COPYRIGHT 2009 ACS on STN

ACCESSION NUMBER: 1919:228 CAPLUS

DOCUMENT NUMBER: 13:228

ORIGINAL REFERENCE NO.: 13:43b-e

TITLE: Transformation of tetrahydronaphthalene (tetralin) in the animal body

AUTHOR(S): Schroeter, G.; Thomas, K.

SOURCE: Z. physiol. Chem. (1918), 101, 262-75

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

AB Tetrahydronaphthalene fed to a dog is absorbed and eventually excreted in combination with urea. The existence of 4 compds. of tetrahydronaphthalene and urea is theoretically possible, and the authors have succeeded in preparing all of them by the interaction of KCNO on the respective amines in H₂O. ar-Tetrahydro- α -carbamidonaphthalene, C₁₁H₁₄ON₂, crystallizes in square plates from alc., soften at 198° and melts at about 206° (quickly heated, at 212°). ar-Tetrahydro- β -carbamidonaphthalene, needles, m. 134° (decomposition). ac-Tetrahydro- α -carbamidonaphthalene, needles, m. 210.5°. ac-Tetrahydro- β -carbamidonaphthalene, needles, m. 183°. Comparison of the natural product with these compds. shows that tetralin in its passage through the body is converted into dl-ac-tetrahydro- α -carbamidonaphthalene. In the preparation of the ar- β -compds., a small amount of a substance was obtained in the form of needles, which did not melt below 245° and possessed the composition of di-ar-tetrahydro- β -naphthylcarbamide, (C₁₀H₁₂N)₂CO.

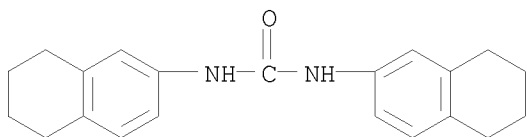
IT 871892-48-7P, Urea, s-bis(5,6,7,8-tetrahydro-2-naphthyl)-

872283-39-1P, Urea, (5,6,7,8-tetrahydro-2-naphthyl)-

RL: PREP (Preparation)
(preparation of)

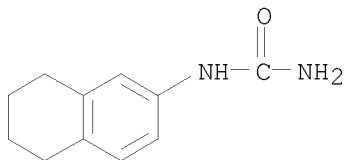
RN 871892-48-7 CAPLUS

CN Urea, N,N'-bis(5,6,7,8-tetrahydro-2-naphthalenyl)- (CA INDEX NAME)



RN 872283-39-1 CAPLUS

CN Urea, N-(5,6,7,8-tetrahydro-2-naphthalenyl)- (CA INDEX NAME)



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FILE 'REGISTRY' ENTERED AT 13:44:55 ON 23 NOV 2009

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L3 4603 S L1 FULL

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L4 116 S L3 FULL
L5 47 S L4 AND PY<2005

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